

(19) World Intellectual Property
Organization
International Bureau



(43) International Publication Date
8 July 2004 (08.07.2004)

PCT

(10) International Publication Number
WO 2004/056829 A1

(51) International Patent Classification⁷: **C07D 487/04**,
A01N 43/90

(21) International Application Number:
PCT/GB2003/005261

(22) International Filing Date: 3 December 2003 (03.12.2003)

(25) Filing Language: English

(26) Publication Language: English

(30) Priority Data:
0230021.8 23 December 2002 (23.12.2002) GB

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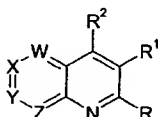
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(81) Designated States (national): AE, AG, AL, AM, AT, AU,
AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR,
CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD,
GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR,
KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN,
MW, MX, MZ, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU,
SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA,
UG, US, UZ, VC, VN, YU, ZA, ZM, ZW.

(84) Designated States (regional): ARIPO patent (BW, GH,
GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW),
Eurasian patent (AM, AZ, BY, KG, KZ, MD, RU, TJ, TM),
European patent (AT, BE, BG, CH, CY, CZ, DE, DK, EE,

[Continued on next page]

(54) Title: FUNGICIDES BASED ON NITROGEN-CONTAINING HETEROCYCLES



(1)

(57) Abstract: Fungicidal compounds having the general formula (1): formula (1) wherein W, Z and one of X and Y are N and the other one of X and Y is CR⁸; R⁸ is H, halo, C₁₋₄ alkyl, C₁₋₄ alkoxy, C₁₋₄ alkylthio or halo(C₁₋₄)alkyl; R and R² are independently H, halo, C₁₋₈ alkyl, C₁₋₈ alkoxy, C₁₋₈ alkylthio, C₂₋₈ alkenyl, C₂₋₈ alkynyl, cyano or NR³R⁴, provided that at least one of R and R² is NR³R⁴; R¹ is halo, C₁₋₈ alkyl, C₂₋₈ alkenyl, C₂₋₈ alkynyl, C₃₋₈ cycloalkyl, C₃₋₈ cycloalkyl(C₁₋₆)alkyl, C₁₋₈ alkoxy, C₁₋₈ alkylthio, aryl, aryloxy, arylthio, heteroaryl, heteroaryloxy, heteroarylthio, aryl(C₁₋₄)alkyl, aryl(C₁₋₄)alkoxy, heteroaryl(C₁₋₄)alkyl, heteroaryl(C₁₋₄)alkoxy, aryl(C₁₋₄)alkylthio, heteroaryl(C₁₋₄)alkylthio, morpholino, piperidino or pyrrolidino; R³ and R⁴ are independently H, C₁₋₈ alkyl, C₂₋₈ alkenyl, C₂₋₈ alkynyl, aryl, aryl(C₁₋₈)alkyl, C₃₋₈ cycloalkyl, C₃₋₈ cycloalkyl(C₁₋₆)alkyl, heteroaryl, heteroaryl(C₁₋₈)alkyl, NR⁵R⁶, provided that not both R³ and R⁴ are H or NR⁵R⁶, or R³ and R⁴ together form a C₃₋₇ alkylene or C₃₋₇ alkenylene chain optionally substituted with one or more C₁₋₄ alkyl or C₁₋₄ alkoxy groups, or, together with the nitrogen atom to which they are attached, R³ and R⁴ form a morpholine, thiomorpholine, thiomorpholine S-oxide or thiomorpholine S-dioxide ring or a piperazine or piperazine N-(C₁₋₄)alkyl (especially N-methyl) ring; and R⁵ and R⁶ are independently H, C₁₋₈ alkyl, C₂₋₈ alkenyl, C₂₋₈ alkynyl, aryl, aryl(C₁₋₈)alkyl, C₃₋₈ cycloalkyl, C₃₋₈ cycloalkyl(C₁₋₆)alkyl, heteroaryl or heteroaryl(C₁₋₈)alkyl; any of the foregoing alkyl, alkenyl, alkynyl or cycloalkyl groups or moieties (other than for R⁸) being optionally substituted with halogen, cyano, C₁₋₆ alkoxy, C₁₋₆ alkylcarbonyl, C₁₋₆ alkoxy carbonyl, C₁₋₆ haloalkoxy, C₁₋₆ alkylthio, tri(C₁₋₄)alkylsilyl, C₁₋₆ alkylamino or C₁₋₆ dialkylamino, any of the foregoing morpholine, thiomorpholine, piperidine, piperazine and pyrrolidine rings being optionally substituted with C₁₋₄ alkyl (especially methyl), and any of the foregoing aryl or heteroaryl groups or moieties being optionally substituted with one or more substituents selected from halo, hydroxy, mercapto, C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₁₋₆ alkoxy, C₂₋₆ alkenyloxy, C₂₋₆ alkynyloxy, halo(C₁₋₆)alkyl, halo(C₁₋₆)alkoxy, C₁₋₆ alkylthio, halo(C₁₋₆)alkylthio, hydroxy(C₁₋₆)alkyl, C₁₋₄ alkoxy(C₁₋₆)alkyl, C₃₋₆ cycloalkyl, C₃₋₆ cycloalkyl(C₁₋₄)alkyl, phenoxy, benzyloxy, benzoyloxy, cyano, isocyanato, thiocyanato, isothiocyanato, nitro, -NR⁷R⁸, -NHCOR⁷, -NHCONR⁷R⁸, -CONR⁷R⁸, -SO₂R⁷, -OSO₂R⁷, -COR⁷, -CR⁷=NR⁸ or -N=CR⁷R⁸, in which R⁷ and R⁸ are independently hydrogen, C₁₋₄ alkyl, halo(C₁₋₄)alkyl, C₁₋₄ alkoxy, halo(C₁₋₄)alkoxy, C₁₋₄ alkylthio, C₃₋₆ cycloalkyl, C₃₋₆ cycloalkyl(C₁₋₄)alkyl, phenyl or benzyl, the phenyl and benzyl groups being optionally substituted with halogen, C₁₋₄ alkyl or C₁₋₄ alkoxy.



ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE,
SI, SK, TR), OAPI patent (BF, BJ, CF, CG, CI, CM, GA,
GN, GQ, GW, ML, MR, NE, SN, TD, TG).

For two-letter codes and other abbreviations, refer to the "Guidance Notes on Codes and Abbreviations" appearing at the beginning of each regular issue of the PCT Gazette.

Published:

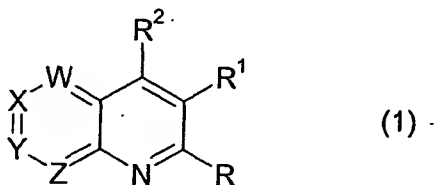
— *with international search report*

FUNGICIDES BASED ON NITROGEN-CONTAINING HETEROCYCLES

This invention relates to novel derivatives of pyridotriazines, to processes for preparing them, to certain intermediate chemicals used in their manufacture, to compositions containing them and to methods of using them to combat fungi, especially fungal infections of plants.

Derivatives of the nitrogen-containing 5,6 ring system *s*-1,2,4-triazolo[1,5-*a*]pyrimidine are known from the patent literature as being useful for controlling phytopathogenic fungi. Examples of recent patent publications include EP-A-1249452, WO 02/051845, WO 02/083676, WO 02/083677, WO 02/088125, WO 02/088126, WO 02/088127.

The present invention is concerned with the provision of novel pyridotriazines for combating phytopathogenic diseases on plants and harvested food crops. Thus, according to the present invention, there is provided a compound of the general formula (1):



wherein

W, Z and one of X and Y are N and the other one of X and Y is CR⁸;

R⁸ is H, halo, C₁₋₄ alkyl, C₁₋₄ alkoxy, C₁₋₄ alkylthio or halo(C₁₋₄)alkyl;

R and R² are independently H, halo, C₁₋₈ alkyl, C₁₋₈ alkoxy, C₁₋₈ alkylthio, C₂₋₈ alkenyl, C₂₋₈ alkynyl, cyano or NR³R⁴, provided that at least one of R and R² is NR³R⁴;

R¹ is halo, C₁₋₈ alkyl, C₂₋₈ alkenyl, C₂₋₈ alkynyl, C₃₋₈ cycloalkyl, C₃₋₈ cycloalkyl(C₁₋₆)alkyl, C₁₋₈ alkoxy, C₁₋₈ alkylthio, aryl, aryloxy, arylthio, heteroaryl, heteroaryloxy, heteroarylthio, aryl(C₁₋₄)alkyl, aryl(C₁₋₄)alkoxy, heteroaryl(C₁₋₄)alkyl, heteroaryl(C₁₋₄)alkoxy, aryl(C₁₋₄)alkylthio, heteroaryl(C₁₋₄)alkylthio, morpholino, piperidino or pyrrolidino;

R³ and R⁴ are independently H, C₁₋₈ alkyl, C₂₋₈ alkenyl, C₂₋₈ alkynyl, aryl, aryl(C₁₋₈)alkyl, C₃₋₈ cycloalkyl, C₃₋₈ cycloalkyl(C₁₋₆)alkyl, heteroaryl, heteroaryl(C₁₋₈)alkyl, NR⁵R⁶, provided that not both R³ and R⁴ are H or NR⁵R⁶, or

R³ and R⁴ together form a C₃₋₇ alkylene or C₃₋₇ alkenylene chain optionally substituted with one or more C₁₋₄ alkyl or C₁₋₄ alkoxy groups, or,

together with the nitrogen atom to which they are attached, R^3 and R^4 form a morpholine, thiomorpholine, thiomorpholine S-oxide or thiomorpholine S-dioxide ring or a piperazine or piperazine *N*-(C_{1-4})alkyl (especially *N*-methyl) ring; and

R^5 and R^6 are independently H, C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, aryl, aryl(C_{1-8})alkyl, C_{3-8} cycloalkyl, C_{3-8} cycloalkyl(C_{1-6})alkyl, heteroaryl or heteroaryl(C_{1-8})alkyl;

any of the foregoing alkyl, alkenyl, alkynyl or cycloalkyl groups or moieties (other than for R^8) being optionally substituted with halogen, cyano, C_{1-6} alkoxy, C_{1-6} alkylcarbonyl, C_{1-6} alkoxycarbonyl, C_{1-6} haloalkoxy, C_{1-6} alkylthio, tri(C_{1-4})alkylsilyl, C_{1-6} alkylamino or C_{1-6} dialkylamino,

any of the foregoing morpholine, thiomorpholine, piperidine, piperazine and pyrrolidine rings being optionally substituted with C_{1-4} alkyl (especially methyl), and

any of the foregoing aryl or heteroaryl groups or moieties being optionally substituted with one or more substituents selected from halo, hydroxy, mercapto, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{1-6} alkoxy, C_{2-6} alkenyloxy, C_{2-6} alkynyloxy, halo(C_{1-6})alkyl, halo(C_{1-6})alkoxy, C_{1-6} alkylthio, halo(C_{1-6})alkylthio, hydroxy(C_{1-6})alkyl, C_{1-4} alkoxy(C_{1-6})alkyl, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl(C_{1-4})alkyl, phenoxy, benzyloxy, benzoyloxy, cyano, isocyano, thiocyanato, iso-

thiocyanato, nitro, $-NR''''R''''$, $-NHCOR''''$, $-NHCONR''''R''''$, $-CONR''''R''''$, $-SO_2R''''$, $-OSO_2R''''$, $-COR''''$, $-CR''''=NR''''$ or $-N=CR''''R''''$, in which R'''' and R'''' are independently hydrogen, C_{1-4} alkyl, halo(C_{1-4})alkyl, C_{1-4} alkoxy, halo(C_{1-4})alkoxy, C_{1-4} alkylthio, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl(C_{1-4})alkyl, phenyl or benzyl, the phenyl and benzyl groups being optionally substituted with halogen, C_{1-4} alkyl or C_{1-4} alkoxy.

The invention includes a compound of the general formula (1) as defined immediately above except that: C_{1-8} alkoxy and C_{1-8} alkylthio are excluded as values of R and R^2 ; C_7 alkylene and C_{3-7} alkenylene are excluded as chains formed by R^3 and R^4 ; the C_{3-6} chain that R^3 and R^4 may form may only be optionally substituted with one or more methyl groups; thiomorpholine, thiomorpholine S-oxide, thiomorpholine S-dioxide and piperazine are excluded as rings that R^3 and R^4 may form; tri(C_{1-4})alkylsilyl is excluded as a substituent of any alkyl, alkenyl, alkynyl or cycloalkyl group or moiety; any morpholine, piperidine or pyrrolidine ring is unsubstituted and C_{1-4} alkylthio is excluded as a value of R^8 .

The compounds of the invention may contain one or more asymmetric carbon atoms and may exist as enantiomers (or as pairs of diastereoisomers) or as mixtures of such. They may also exist as diastereoisomers by virtue of restricted rotation about a bond. However,

mixtures of enantiomers or diastereoisomers may be separated into individual isomers or isomer pairs, and this invention embraces such isomers and mixtures thereof in all proportions. It is to be expected that for any given compound, one isomer may be more fungicidally active than another.

5 Except where otherwise stated, alkyl groups and alkyl moieties of alkoxy, alkylthio, etc., contain from 1 to 8, suitably from 1 to 6 and typically from 1 to 4, carbon atoms in the form of straight or branched chains. Examples are methyl, ethyl, *n*- and *iso*-propyl, *n*-, *sec*-, *iso*- and *tert*-butyl, *n*-pentyl and *n*-hexyl. Cycloalkyl groups contain from 3 to 8, typically from 3 to 6, carbon atoms and include bicycloalkyl groups such as the bicyclo[2.2.1]heptyl
10 group. Haloalkyl groups or moieties are typically trichloromethyl or trifluoromethyl or contain a trichloromethyl or trifluoromethyl terminal group.

Except where otherwise stated, alkenyl and alkynyl moieties also contain from 2 to 8, suitably from 2 to 6 and typically from 2 to 4, carbon atoms in the form of straight or branched chains. Examples are allyl, 2-methylallyl and propargyl. Optional substituents
15 include halo, typically fluoro. An example of halo-substituted alkenyl is 3,4,4-trifluoro-*n*-butenyl.

Halo includes fluoro, chloro, bromo and iodo. Most commonly it is fluoro, chloro or bromo and usually fluoro or chloro.

Aryl is usually phenyl but also includes naphthyl, anthryl and phenanthryl.

20 Heteroaryl is typically a 5- or 6-membered aromatic ring containing one or more O, N or S heteroatoms, which may be fused to one or more other aromatic or heteroaromatic rings, such as a benzene ring. Examples are thienyl, furyl, pyrrolyl, isoxazolyl, oxazolyl, oxadiazolyl, pyrazolyl, imidazolyl, triazolyl, isothiazolyl, tetrazolyl, thiadiazolyl, pyridyl, pyrimidinyl, pyrazinyl, pyridazinyl, triazinyl, benzofuryl, benzothienyl, dibenzofuryl,
25 benzothiazolyl, benzoxazolyl, benzimidazolyl, indolyl, quinoliny and quinoxaliny groups and, where appropriate, *N*-oxides thereof.

The 6,6-ring systems embraced by the general formula (1) are pyrido[2,3-*e*][1,2,4]-triazines (where Y is CR⁸ and W, X and Z are all N) and pyrido[3,2-*e*][1,2,4]triazines (where X is CR⁸ and W, Y and Z are all N).

30 R⁸ is H, halo (for example bromo), C₁₋₄ alkyl (for example methyl), C₁₋₄ alkoxy (for example methoxy), C₁₋₄ alkylthio (for example methylthio) or halo(C₁₋₄)alkyl (for example trifluoromethyl). Usually R⁸ will be H.

One of R and R^2 , preferably R^2 , is NR^3R^4 . The other is typically halo, especially chloro or fluoro. The more active compounds are usually those where R^2 is NR^3R^4 . R^3 is typically C_{1-8} alkyl (for example ethyl, *n*-propyl, *n*-butyl, *sec*-butyl (the *S*- or *R*-isomer or the racemate) and *tert*-butyl), halo(C_{1-8})alkyl (for example 2,2,2-trifluoroethyl, 2,2,2-trifluoro-1-methylethyl (the *S*- or *R*-isomer or the racemate), 3,3,3-trifluoropropyl and 4,4,4-trifluorobutyl), hydroxy(C_{1-8})alkyl (for example hydroxyethyl), C_{1-4} alkoxy(C_{1-8})alkyl (for example methoxymethyl and methoxy-*iso*-butyl), C_{1-4} alkoxyhalo(C_{1-8})alkyl (for example 2-methoxy-2-trifluoromethylethyl), tri(C_{1-4})alkylsilyl(C_{1-6})alkyl (for example trimethylsilylmethyl), C_{1-4} alkylcarbonyl(C_{1-8})alkyl (for example 1-acetyethyl and 1-*tert*-butylcarbonylethyl), C_{1-4} alkylcarbonylhalo(C_{1-8})alkyl (for example 1-acetyl-2,2,2-trifluoroethyl), phenyl($_{1-4}$)alkyl (for example benzyl), C_{2-8} alkenyl (for example allyl and methylallyl), halo(C_{2-8})alkenyl (for example 3-methyl-4,4-difluorobut-3-enyl), C_{2-8} alkynyl (for example propargyl), C_{3-8} cycloalkyl (for example cyclopropyl, cyclobutyl, cyclopentyl and cyclohexyl) optionally substituted with chloro, fluoro or methyl, C_{3-8} cycloalkyl(C_{1-4})alkyl (for example cyclopropylmethyl, cyclobutylmethyl, cyclopentylmethyl and cyclohexylmethyl), phenylamino, piperidino or morpholino, the phenyl ring of phenylalkyl or phenylamino being optionally substituted with one, two or three substituents selected from halo (typically fluoro, chloro or bromo), C_{1-4} alkyl (typically methyl), halo(C_{1-4})alkyl (typically trifluoromethyl), C_{1-4} alkoxy (typically methoxy) and halo(C_{1-4})alkoxy (typically trifluoromethoxy). R^4 is typically H, C_{1-4} alkyl (for example ethyl and *n*-propyl), halo(C_{1-4})alkyl (for example 2,2,2-trifluoroethyl) or amino. Alternatively R^3 and R^4 together form a C_{4-6} alkylene chain optionally substituted with methyl, for example 3-methylpentylene, or, together with the nitrogen atom to which they are attached, R^3 and R^4 form a morpholine, thiomorpholine, thiomorpholine *S*-oxide or thiomorpholine *S*-dioxide ring or a piperazine or piperazine *N*-(C_{1-4})alkyl (especially *N*-methyl) ring, in which the morpholine or piperazine rings are optionally substituted with methyl.

Typically R^1 is phenyl optionally substituted with from one to five halogen atoms, particularly fluorine and chlorine atoms and especially fluorine atoms or with from one to three substituents selected from halo (for example fluoro and chloro), C_{1-4} alkyl (for example methyl), halo(C_{1-4})alkyl (for example trifluoromethyl), C_{1-4} alkoxy (for example methoxy) or halo(C_{1-4})alkoxy (for example trifluoromethoxy). Examples are 2,6-difluorophenyl, 2-fluoro-6-chlorophenyl, 2,5,6-trifluorophenyl, 2,4,6-trifluorophenyl, 2,6-difluoro-4-methoxyphenyl,

pentafluorophenyl, 2-fluorophenyl, 2,3,5,6-tetrafluorophenyl, 2-chloro-4,6-difluorophenyl, 2-chlorophenyl, 2,6-dichlorophenyl, 2,4-dichlorophenyl, 2,4,6-trichlorophenyl, 2,3,6-trichlorophenyl, pentachlorophenyl, 2-fluoro-4,6-dichlorophenyl, 4-fluoro-2,6-dichlorophenyl, 2-bromophenyl, 2-fluoro-6-bromophenyl, 2-bromo-4,6-difluorophenyl, 2-fluoro-6-methylphenyl, 2-chloro-6-methylphenyl, 2-methoxyphenyl, 2,6-dimethoxyphenyl, 2-fluoro-6-methoxyphenyl, 2-trifluoromethylphenyl, 2-fluoro-6-trifluoromethylphenyl, 2,6-di-(trifluoromethyl)phenyl, 2-chloro-6-trifluoromethylphenyl, 2,4-difluoro-6-trifluoromethylphenyl, 2,4-difluoro-6-methoxyphenyl and 2,4-difluoro-6-methylphenyl.

Also of particular interest are compounds where R^1 is pyridyl optionally substituted with from one to four halogen atoms or with from one to three substituents selected from halo (for example fluoro and chloro), C_{1-4} alkyl (for example methyl), halo(C_{1-4})alkyl (for example trifluoromethyl), C_{1-4} alkoxy (for example methoxy) or halo(C_{1-4})alkoxy (for example trifluoromethoxy). Examples are 2,4-difluoropyrid-3-yl, 3,5-difluoropyrid-4-yl, tetrafluoropyrid-4-yl, 3-fluoropyrid-2-yl, 4-fluoropyrid-3-yl, 3-fluoropyrid-4-yl, 2-fluoropyrid-3-yl, 2,4,6-trifluoropyrid-3-yl, 3,5-difluoropyrid-2-yl, 2,6-difluoropyrid-3-yl, 2,4-difluoro-6-methoxypyrid-3-yl, 2-fluoro-4-chloropyrid-3-yl, 3-fluoro-5-chloropyrid-4-yl, 2-chloro-4-fluoropyrid-3-yl, 2,4-dichloropyrid-3-yl, 3-chloropyrid-2-yl, 4-chloropyrid-3-yl, 3-chloropyrid-4-yl, 2-chloropyrid-3-yl, 3-trifluoromethylpyrid-2-yl, 4-trifluoromethylpyrid-3-yl, 3,5-dichloropyrid-2-yl, 4,6-dichloropyrid-3-yl, 3-trifluoromethylpyrid-4-yl, 2-trifluoromethylpyrid-3-yl, 2-fluoro-4-trifluoromethylpyrid-3-yl, 3-fluoro-5-trifluoromethylpyrid-4-yl, 4-fluoro-2-trifluoromethylpyrid-3-yl, 2,6-dichloropyrid-3-yl, 3,5-dichloropyrid-4-yl, 3-chloro-6-trifluoromethylpyrid-2-yl, 3-fluoro-6-trifluoromethylpyrid-2-yl, pyrid-2-yl, pyrid-3-yl and pyrid-4-yl.

Also of particular interest are compounds where R^1 is 2- or 3-thienyl optionally substituted with from one to three halogen atoms or with from one to three substituents selected from halo (for example fluoro and chloro), C_{1-4} alkyl (for example methyl), halo(C_{1-4})alkyl (for example trifluoromethyl), C_{1-4} alkoxy (for example methoxy) or halo(C_{1-4})alkoxy (for example trifluoromethoxy). Examples are 3-fluorothien-2-yl, 3-chlorothien-2-yl, 2,4-difluorothien-3-yl, 2,4-dichlorothien-3-yl and 2,4,5-trichlorothien-3-yl.

Examples of other values of R^1 of especial interest are unsubstituted piperidino and morpholino, 2-methylpiperidino, 2,6-dimethylpiperidino and 2,6-dimethylmorpholino.

In one aspect the invention provides a compound of the general formula (1) wherein W, Z and one of X and Y are N and the other one of X and Y is CR⁸;
R⁸ is H, halo, C₁₋₄ alkyl, C₁₋₄ alkoxy, C₁₋₄ alkylthio or halo(C₁₋₄)alkyl;
one of R and R² (preferably R²) is NR³R⁴ and the other is halo;
5 R¹ is halo, C₁₋₈ alkyl, C₂₋₈ alkenyl, C₂₋₈ alkynyl, C₃₋₈ cycloalkyl, C₃₋₈ cycloalkyl(C₁₋₆)alkyl, C₁₋₈ alkoxy, C₁₋₈ alkylthio, aryl, aryloxy, arylthio, heteroaryl, heteroaryloxy, heteroarylthio, aryl(C₁₋₄)alkyl, aryl(C₁₋₄)alkoxy, heteroaryl(C₁₋₄)alkyl, heteroaryl(C₁₋₄)alkoxy, aryl(C₁₋₄)alkylthio, heteroaryl(C₁₋₄)alkylthio, morpholino, piperidino or pyrrolidino;
R³ and R⁴ are independently H, C₁₋₈ alkyl, C₂₋₈ alkenyl, C₂₋₈ alkynyl, aryl, aryl(C₁₋₈)alkyl, C₃₋₈
10 cycloalkyl, C₃₋₈ cycloalkyl(C₁₋₆)alkyl, heteroaryl, heteroaryl(C₁₋₈)alkyl, NR⁵R⁶, provided that not both R³ and R⁴ are H or NR⁵R⁶, or
R³ and R⁴ together form a C₃₋₇ alkylene or C₃₋₇ alkenylene chain optionally substituted with one or more C₁₋₄ alkyl or C₁₋₄ alkoxy groups, or,
together with the nitrogen atom to which they are attached, R³ and R⁴ form a morpholine,
15 thiomorpholine, thiomorpholine S-oxide or thiomorpholine S-dioxide ring or a piperazine or piperazine N-(C₁₋₄)alkyl (especially N-methyl) ring; and
R⁵ and R⁶ are independently H, C₁₋₈ alkyl, C₂₋₈ alkenyl, C₂₋₈ alkynyl, aryl, aryl(C₁₋₈)alkyl, C₃₋₈ cycloalkyl, C₃₋₈ cycloalkyl(C₁₋₆)alkyl, heteroaryl or heteroaryl(C₁₋₈)alkyl;
any of the foregoing alkyl, alkenyl, alkynyl or cycloalkyl groups or moieties (other than for
20 R⁸) being optionally substituted with halogen, cyano, C₁₋₆ alkoxy, C₁₋₆ alkylcarbonyl, C₁₋₆ alkoxy carbonyl, C₁₋₆ haloalkoxy, C₁₋₆ alkylthio, tri(C₁₋₄)alkylsilyl, C₁₋₆ alkylamino or C₁₋₆ dialkylamino,
any of the foregoing morpholine, thiomorpholine, piperidine, piperazine and pyrrolidine rings being optionally substituted with C₁₋₄ alkyl (especially methyl), and
25 any of the foregoing aryl, heteroaryl, aryloxy or heteroaryl groups being optionally substituted with one or more substituents selected from halo, hydroxy, mercapto, C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₁₋₆ alkoxy, C₂₋₆ alkenyloxy, C₂₋₆ alkynyloxy, halo(C₁₋₆)alkyl, halo(C₁₋₆)alkoxy, C₁₋₆ alkylthio, halo(C₁₋₆)alkylthio, hydroxy(C₁₋₆)alkyl, C₁₋₄ alkoxy(C₁₋₆)alkyl, C₃₋₆ cycloalkyl, C₃₋₆ cycloalkyl(C₁₋₄)alkyl, phenoxy, benzyloxy, benzoyloxy, cyano,
30 isocyano, thiocyanato, isothiocyanato, nitro, -NR^{'''}R^{'''}, -NHCOR^{'''}, -NHCONR^{'''}R^{'''}, -CONR^{'''}R^{'''}, -SO₂R^{'''}, -OSO₂R^{'''}, -COR^{'''}, -CR^{'''}=NR^{'''} or -N=CR^{'''}R^{'''}, in which R^{'''} and R^{'''} are independently hydrogen, C₁₋₄ alkyl, halo(C₁₋₄)alkyl, C₁₋₄ alkoxy, halo(C₁₋₄)alkoxy, C₁₋₄

alkylthio, C₃₋₆ cycloalkyl, C₃₋₆ cycloalkyl(C₁₋₄)alkyl, phenyl or benzyl, the phenyl and benzyl groups being optionally substituted with halogen, C₁₋₄ alkyl or C₁₋₄ alkoxy.

Of particular interest are compounds where R⁸ is H.

The invention includes a compound of the general formula (1) as defined immediately
5 above except that: C₇ alkylene and C₃₋₇ alkenylene are excluded as chains formed by R³ and R⁴; the C₃₋₆ chain that R³ and R⁴ may form may only be optionally substituted with one or more methyl groups; thiomorpholine, thiomorpholine S-oxide, thiomorpholine S-dioxide and piperazine are excluded as rings that R³ and R⁴ may form; tri(C₁₋₄)alkylsilyl is excluded as a substituent of any alkyl, alkenyl, alkynyl or cycloalkyl group or moiety; any morpholine,
10 piperidine or pyrrolidine ring is unsubstituted and C₁₋₄ alkylthio is excluded as a value of R⁸.

In another aspect the invention provides a compound of the general formula (1) wherein

W, Z and one of X and Y are N and the other one of X and Y is CR⁸;

R⁸ is H, halo, C₁₋₄ alkyl, C₁₋₄ alkoxy, C₁₋₄ alkylthio or halo(C₁₋₄)alkyl;

15 one of R and R² (preferably R²) is NR³R⁴ and the other is halo;

R¹ is halo, C₁₋₈ alkyl, C₂₋₈ alkenyl, C₂₋₈ alkynyl, C₃₋₈ cycloalkyl, C₃₋₈ cycloalkyl(C₁₋₆)alkyl, C₁₋₈ alkoxy, C₁₋₈ alkylthio, aryl, aryloxy, arylthio, heteroaryl, heteroaryloxy, heteroarylthio, aryl(C₁₋₄)alkyl, aryl(C₁₋₄)alkoxy, heteroaryl(C₁₋₄)alkyl, heteroaryl(C₁₋₄)alkoxy, aryl(C₁₋₄)alkylthio, heteroaryl(C₁₋₄)alkylthio, morpholino, piperidino or pyrrolidino;

20 R³ is C₁₋₄ alkyl, halo(C₁₋₄)alkyl, C₂₋₄ alkenyl, C₃₋₆ cycloalkyl, C₃₋₆ cycloalkyl(C₁₋₄)alkyl or phenylamino in which the phenyl ring is optionally substituted with one, two or three substituents selected from halo, C₁₋₄ alkyl, halo(C₁₋₄)alkyl, C₁₋₄ alkoxy and halo(C₁₋₄)alkoxy; and R⁴ is H, C₁₋₄ alkyl or amino, or

R³ and R⁴ together form a C₄₋₆ alkylene chain optionally substituted with C₁₋₄ alkyl or C₁₋₄
25 alkoxy, or,

together with the nitrogen atom to which they are attached, R³ and R⁴ form a morpholine, thiomorpholine, thiomorpholine S-oxide or thiomorpholine S-dioxide ring or a piperazine or piperazine N-(C₁₋₄)alkyl (especially N-methyl) ring;

any of the foregoing alkyl, alkenyl, alkynyl or cycloalkyl groups or moieties (other than for
30 R⁸) being optionally substituted with halogen, cyano, C₁₋₆ alkoxy, C₁₋₆ alkylcarbonyl, C₁₋₆ alkoxy carbonyl, C₁₋₆ haloalkoxy, C₁₋₆ alkylthio, tri(C₁₋₄)alkylsilyl, C₁₋₆ alkylamino or C₁₋₆ dialkylamino,

any of the foregoing morpholine, thiomorpholine, piperidine, piperazine and pyrrolidine rings being optionally substituted with C₁₋₄ alkyl (especially methyl), and
 any of the foregoing aryl or heteroaryl groups or moieties being optionally substituted with one or more substituents selected from halo, hydroxy, mercapto, C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₁₋₆ alkoxy, C₂₋₆ alkenyloxy, C₂₋₆ alkynyloxy, halo(C₁₋₆)alkyl, halo(C₁₋₆)alkoxy, C₁₋₆ alkylthio, halo(C₁₋₆)alkylthio, hydroxy(C₁₋₆)alkyl, C₁₋₄ alkoxy(C₁₋₆)alkyl, C₃₋₆ cycloalkyl, C₃₋₆ cycloalkyl(C₁₋₄)alkyl, phenoxy, benzyloxy, benzoyloxy, cyano, isocyano, thiocyanato, isothiocyanato, nitro, -NR^{'''}R^{'''}, -NHCOR^{'''}, -NHCONR^{'''}R^{'''}, -CONR^{'''}R^{'''}, -SO₂R^{'''}, -OSO₂R^{'''}, -COR^{'''}, -CR^{'''}=NR^{'''} or -N=CR^{'''}R^{'''}, in which R^{'''} and R^{'''} are independently hydrogen, C₁₋₄ alkyl, halo(C₁₋₄)alkyl, C₁₋₄ alkoxy, halo(C₁₋₄)alkoxy, C₁₋₄ alkylthio, C₃₋₆ cycloalkyl, C₃₋₆ cycloalkyl(C₁₋₄)alkyl, phenyl or benzyl, the phenyl and benzyl groups being optionally substituted with halogen, C₁₋₄ alkyl or C₁₋₄ alkoxy.

Of particular interest are compounds where R⁸ is H.

The invention includes a compound of the general formula (1) as defined immediately above except that: the C₄₋₆ chain that R³ and R⁴ may form may only be optionally substituted with methyl; thiomorpholine, thiomorpholine S-oxide, thiomorpholine S-dioxide and piperazine are excluded as rings that R³ and R⁴ may form; tri(C₁₋₄)alkylsilyl is excluded as a substituent of any alkyl, alkenyl, alkynyl or cycloalkyl group or moiety; any morpholine, piperidine or pyrrolidine ring is unsubstituted and C₁₋₄ alkylthio is excluded as a value of R⁸.

In yet another aspect the invention provides a compound of the general formula (1) wherein

W, Z and one of X and Y are N and the other one of X and Y is CR⁸;

R⁸ is H, halo, C₁₋₄ alkyl, C₁₋₄ alkoxy, C₁₋₄ alkylthio or halo(C₁₋₄)alkyl;

R and R² are independently H, halo, C₁₋₈ alkyl, C₁₋₈ alkoxy, C₁₋₈ alkylthio, C₂₋₈ alkenyl, C₂₋₈ alkynyl, cyano or NR³R⁴, provided that at least one of R and R² (preferably R²) is NR³R⁴;

R¹ is optionally substituted phenyl;

R³ and R⁴ are independently H, C₁₋₈ alkyl, C₂₋₈ alkenyl, C₂₋₈ alkynyl, aryl, aryl(C₁₋₈)alkyl, C₃₋₈ cycloalkyl, C₃₋₈ cycloalkyl(C₁₋₆)alkyl, heteroaryl, heteroaryl(C₁₋₈)alkyl, NR⁵R⁶, provided that not both R³ and R⁴ are H or NR⁵R⁶, or

R³ and R⁴ together form a C₃₋₇ alkylene or C₃₋₇ alkenylene chain optionally substituted with one or more C₁₋₄ alkyl or C₁₋₄ alkoxy groups, or, together with the nitrogen atom to which they are attached, R³ and R⁴ form a morpholine,

thiomorpholine, thiomorpholine S-oxide or thiomorpholine S-dioxide ring or a piperazine or piperazine *N*-(C₁₋₄)alkyl (especially *N*-methyl) ring; and

R⁵ and R⁶ are independently H, C₁₋₈ alkyl, C₂₋₈ alkenyl, C₂₋₈ alkynyl, aryl, aryl(C₁₋₈)alkyl, C₃₋₈ cycloalkyl, C₃₋₈ cycloalkyl(C₁₋₆)alkyl, heteroaryl or heteroaryl(C₁₋₈)alkyl;

5 any of the foregoing alkyl, alkenyl, alkynyl or cycloalkyl groups or moieties (other than for R⁸) being optionally substituted with halogen, cyano, C₁₋₆ alkoxy, C₁₋₆ alkylcarbonyl, C₁₋₆ alkoxycarbonyl, C₁₋₆ haloalkoxy, C₁₋₆ alkylthio, tri(C₁₋₄)alkylsilyl, C₁₋₆ alkylamino or C₁₋₆ dialkylamino,

any of the foregoing morpholine, thiomorpholine, piperidine, piperazine and pyrrolidine rings
10 being optionally substituted with C₁₋₄ alkyl (especially methyl), and

any of the foregoing aryl or heteroaryl groups or moieties, including the phenyl group of R¹, being optionally substituted with one or more substituents selected from halo, hydroxy, mercapto, C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₁₋₆ alkoxy, C₂₋₆ alkenyloxy, C₂₋₆ alkynyloxy, halo(C₁₋₆)alkyl, halo(C₁₋₆)alkoxy, C₁₋₆ alkylthio, halo(C₁₋₆)alkylthio, hydroxy(C₁₋₆)alkyl, C₁₋₄
15 alkoxy(C₁₋₆)alkyl, C₃₋₆ cycloalkyl, C₃₋₆ cycloalkyl(C₁₋₄)alkyl, phenoxy, benzyloxy, benzoyloxy, cyano, isocyano, thiocyanato, isothiocyanato, nitro, -NR^{'''}R^{'''}, -NHCOR^{'''}, -NHCONR^{'''}R^{'''}, -CONR^{'''}R^{'''}, -SO₂R^{'''}, -OSO₂R^{'''}, -COR^{'''}, -CR^{'''}=NR^{'''} or -N=CR^{'''}R^{'''}, in which R^{'''} and R^{'''} are independently hydrogen, C₁₋₄ alkyl, halo(C₁₋₄)alkyl, C₁₋₄ alkoxy, halo(C₁₋₄)alkoxy, C₁₋₄ alkylthio, C₃₋₆ cycloalkyl, C₃₋₆ cycloalkyl(C₁₋₄)alkyl, phenyl or benzyl, the
20 phenyl and benzyl groups being optionally substituted with halogen, C₁₋₄ alkyl or C₁₋₄ alkoxy. Of particular interest are compounds where R⁸ is H.

The invention includes a compound of the general formula (1) as defined immediately above except that: C₁₋₈ alkoxy and C₁₋₈ alkylthio are excluded as values of R and R²; C₇ alkylene and C₃₋₇ alkenylene are excluded as chains formed by R³ and R⁴; the C₃₋₆ chain that
25 R³ and R⁴ may form may only be optionally substituted with one or more methyl groups; thiomorpholine, thiomorpholine S-oxide, thiomorpholine S-dioxide and piperazine are excluded as rings that R³ and R⁴ may form; tri(C₁₋₄)alkylsilyl is excluded as a substituent of any alkyl, alkenyl, alkynyl or cycloalkyl group or moiety; the morpholine ring that R³ and R⁴ may form is unsubstituted and C₁₋₄ alkylthio is excluded as a value of R⁸.

30 In still yet another aspect the invention provides a compound of the general formula (1) wherein

W, Z and one of X and Y are N and the other one of X and Y is CR⁸;

R^8 is H, halo (e.g. fluoro, chloro or bromo), C_{1-4} alkyl (e.g. methyl), C_{1-4} alkoxy (e.g. methoxy), C_{1-4} alkylthio (e.g. methylthio) or halo(C_{1-4})alkyl (e.g. trifluoromethyl);

R is H, halo (e.g. fluoro, chloro or bromo), C_{1-4} alkyl (e.g. methyl), C_{1-4} alkoxy (e.g. methoxy) or cyano;

- 5 R^1 is phenyl optionally substituted with from one to five halogen atoms or with from one to three substituents selected from halo, C_{1-4} alkyl, halo(C_{1-4})alkyl, C_{1-4} alkoxy or halo(C_{1-4})-alkoxy, pyridyl optionally substituted with from one to four halogen atoms or with from one to three substituents selected from halo, C_{1-4} alkyl, halo(C_{1-4})alkyl, C_{1-4} alkoxy or halo(C_{1-4})-alkoxy, 2- or 3-thienyl optionally substituted with from one to three halogen atoms or with
10 from one to three substituents selected from halo, C_{1-4} alkyl, halo(C_{1-4})alkyl, C_{1-4} alkoxy or halo(C_{1-4})alkoxy, or piperidino or morpholino both optionally substituted with one or two methyl groups;

R^2 is NR^3R^4 ;

- R^3 is C_{1-8} alkyl, halo(C_{1-8})alkyl, hydroxy(C_{1-8})alkyl, C_{1-4} alkoxy(C_{1-8})alkyl, C_{1-4} alkoxyhalo-
15 (C_{1-8})alkyl, tri(C_{1-4})alkylsilyl(C_{1-6})alkyl, C_{1-4} alkylcarbonyl(C_{1-8})alkyl, C_{1-4} alkylcarbonylhalo(C_{1-8})alkyl, phenyl($_{1-4}$)alkyl, C_{2-8} alkenyl, halo(C_{2-8})alkenyl, C_{2-8} alkynyl, C_{3-8} cycloalkyl optionally substituted with chloro, fluoro or methyl, C_{3-8} cycloalkyl(C_{1-4})alkyl, phenylamino, piperidino or morpholino, the phenyl ring of phenylalkyl or phenylamino being optionally substituted with one, two or three substituents selected from halo, C_{1-4} alkyl, halo(C_{1-4})alkyl,
20 C_{1-4} alkoxy and halo(C_{1-4})alkoxy; and

R^4 is H, C_{1-4} alkyl, halo(C_{1-4})alkyl or amino, or

R^3 and R^4 together form a C_{3-7} alkylene or C_{3-7} alkenylene chain optionally substituted with methyl; or,

- together with the nitrogen atom to which they are attached, R^3 and R^4 form a morpholine,
25 thiomorpholine, thiomorpholine S-oxide or thiomorpholine S-dioxide ring or a piperazine or piperazine *N*-(C_{1-4})alkyl (especially *N*-methyl) ring, in which the morpholine or piperazine rings are optionally substituted with methyl.

Of particular interest are compounds where R^8 is H.

- In still yet another aspect the invention provides a compound of the general formula
30 (1) wherein

W, Z and one of X and Y are N and the other one of X and Y is CR^8 ;

R^8 is H, halo, C_{1-4} alkyl, C_{1-4} alkoxy or halo(C_{1-4})alkyl;

R is halo;

R¹ is phenyl optionally substituted with from one to five halogen atoms or with from one to three substituents selected from halo, C₁₋₄ alkyl, halo(C₁₋₄)alkyl, C₁₋₄ alkoxy or halo(C₁₋₄)alkoxy;

5 R² is NR³R⁴;

R³ is C₁₋₄ alkyl, halo(C₁₋₄)alkyl, C₂₋₄ alkenyl, C₃₋₆ cycloalkyl, C₃₋₆ cycloalkyl(C₁₋₄)alkyl or phenylamino in which the phenyl ring is optionally substituted with one, two or three substituents selected from halo, C₁₋₄ alkyl, halo(C₁₋₄)alkyl, C₁₋₄ alkoxy and halo(C₁₋₄)alkoxy; and

10 R⁴ is H, C₁₋₄ alkyl or amino, or R³ and R⁴ together form a C₄₋₆ alkylene chain optionally substituted with methyl, or, together with the nitrogen atom to which they are attached, R³ and R⁴ form a morpholine ring.

Of particular interest are compounds where R⁸ is H.

Compounds that form part of the invention are illustrated in Tables 1 to 130 below.

15 Characterising data are given later in the Examples and in Table 136.

In Table 1 the compounds have the general formula (1A), where W, X and Z are N and Y is CH, R is Cl, R¹ is 2,4,6-trifluorophenyl and R³ and R⁴ are as shown in the table.

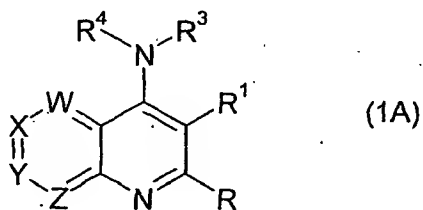


Table 1

| Cmpd No | R ³ | R ⁴ |
|---------|----------------------------------------------------------------------------------------|-----------------------------------------|
| 1 | C ₂ H ₅ | H |
| 2 | <i>n</i> -C ₃ H ₇ | H |
| 3 | <i>i</i> -C ₃ H ₇ | H |
| 4 | <i>n</i> -C ₄ H ₉ | H |
| 5 | <i>i</i> -C ₄ H ₉ | H |
| 6 | CH ₂ =CHCH ₂ | H |
| 7 | CH ₂ =C(CH ₃)CH ₂ | H |
| 8 | CF ₃ CH ₂ | H |
| 9 | CF ₃ CH ₂ CH ₂ | H |
| 10 | CF ₃ CH ₂ CH ₂ CH ₂ | H |
| 11 | CF ₃ (CH ₃)CH | H |
| 12 | (S)-CF ₃ (CH ₃)CH | H |
| 13 | (R)-CF ₃ (CH ₃)CH | H |
| 14 | cyclo-C ₃ H ₅ | H |
| 15 | cyclo-C ₄ H ₇ | H |
| 16 | cyclo-C ₅ H ₉ | H |
| 17 | cyclo-C ₆ H ₁₁ | H |
| 18 | cyclo-C ₃ H ₅ CH ₂ | H |
| 19 | cyclo-C ₄ H ₇ CH ₂ | H |
| 20 | -(CH ₂) ₂ O(CH ₂) ₂ - | |
| 21 | cyclo-C ₆ H ₁₁ CH ₂ | H |
| 22 | -(CH ₂) ₂ CH(CH ₃)(CH ₂) ₂ - | |
| 23 | CH ₃ CH ₂ (CH ₃)CH | H |
| 24 | (S)-CH ₃ CH ₂ (CH ₃)CH | H |
| 25 | (R)-CH ₃ CH ₂ (CH ₃)CH | H |
| 26 | C ₂ H ₅ | C ₂ H ₅ |
| 27 | <i>n</i> -C ₃ H ₇ | <i>n</i> -C ₃ H ₇ |
| 28 | CH ₂ =C(CH ₃)CH ₂ | C ₂ H ₅ |
| 29 | CF ₃ CH ₂ | C ₂ H ₅ |
| 30 | C ₂ H ₅ | NH ₂ |
| 31 | <i>n</i> -C ₃ H ₇ | NH ₂ |
| 32 | <i>i</i> -C ₃ H ₇ | NH ₂ |
| 33 | <i>n</i> -C ₄ H ₉ | NH ₂ |
| 34 | CH ₂ =CHCH ₂ | NH ₂ |
| 35 | CH ₂ =C(CH ₃)CH ₂ | NH ₂ |

| Cmpd No | R ³ | R ⁴ |
|---------|------------------------------------------------------------------------------|-----------------|
| 36 | CF ₂ =CFCH ₂ CH ₂ | NH ₂ |
| 37 | CF ₃ CH ₂ | NH ₂ |
| 38 | CF ₃ CH ₂ CH ₂ | NH ₂ |
| 39 | CF ₃ CH ₂ CH ₂ CH ₂ | NH ₂ |
| 40 | 4- <i>t</i> -C ₄ H ₉ -C ₆ H ₄ NH | H |
| 41 | 4-F-C ₆ H ₄ NH | H |
| 42 | C ₆ H ₅ NH | H |
| 43 | 4-CH ₃ -C ₆ H ₄ NH | H |
| 44 | 4-Br-C ₆ H ₄ NH | H |
| 45 | 2-F-C ₆ H ₄ NH | H |
| 46 | 3,4-Cl ₂ -C ₆ H ₃ NH | H |
| 47 | 3-CF ₃ -C ₆ H ₄ NH | H |
| 48 | 3,5-Cl ₂ -C ₆ H ₃ NH | H |
| 49 | 4-CF ₃ O-C ₆ H ₃ NH | H |
| 50 | 2-CF ₃ -C ₆ H ₄ NH | H |
| 51 | 4-CF ₃ -C ₆ H ₄ NH | H |
| 52 | 2-Br-C ₆ H ₄ NH | H |
| 53 | 2-Cl-C ₆ H ₄ NH | H |
| 54 | 2-CH ₃ -4-Cl-C ₆ H ₃ NH | H |
| 55 | 2-CH ₃ -5-F-C ₆ H ₃ NH | H |
| 56 | 3-Cl-C ₆ H ₄ NH | H |
| 57 | CH ₃ | H |
| 58 | (CH ₃) ₂ CHCH ₂ | H |
| 59 | (CH ₃) ₃ CCH ₂ | H |
| 60 | (CH ₃) ₃ C(CH ₃)CH | H |
| 61 | CH ₃ CH ₂ (CH ₃) ₂ C | H |
| 62 | CH ₃ CH ₂ (CF ₃)CH | H |
| 63 | (S)-CH ₃ CH ₂ (CF ₃)CH | H |
| 64 | (R)-CH ₃ CH ₂ (CF ₃)CH | H |
| 65 | CH ₃ CH ₂ (CH ₃ CH ₂)CH | H |
| 66 | (CH ₃) ₂ CH(CH ₃ CH ₂)CH | H |
| 67 | (CH ₃) ₂ CH(CH ₃)CH | H |
| 68 | (CH ₃) ₂ CH(CF ₃)CH | H |
| 69 | (S)-(CH ₃) ₂ CH(CF ₃)CH | H |
| 70 | (R)-(CH ₃) ₂ CH(CF ₃)CH | H |

| Cmpd No | R ³ | R ⁴ |
|---------|------------------------------------------------------------------------------------------|----------------|
| 71 | HC=C(CH ₃)CH ₂ | H |
| 72 | CH ₂ =CH(CH ₃ CH ₂)CH | H |
| 73 | CH ₃ CH ₂ CH ₂ (CH ₃)CH | H |
| 74 | CH ₃ CH ₂ CH ₂ (CF ₃)CH | H |
| 75 | (S)-CH ₃ CH ₂ CH ₂ (CF ₃)CH | H |
| 76 | (R)-CH ₃ CH ₂ CH ₂ (CF ₃)CH | H |
| 77 | CH ₃ CH ₂ CH ₂ (CH ₃) ₂ C | H |
| 78 | CH ₃ CH ₂ (CH ₃)CHCH ₂ | H |
| 79 | (CH ₃) ₂ CHCH ₂ CH ₂ | H |
| 80 | (CH ₃) ₃ CCH ₂ CH ₂ | H |
| 81 | CH ₃ CH ₂ (CH ₃)CH(CH ₃)-CH | H |
| 82 | CH ₃ CH ₂ (CH ₃)CH(CF ₃)-CH | H |
| 83 | (S)-CH ₃ CH ₂ (CH ₃)CH-(CF ₃)CH | H |
| 84 | (R)-CH ₃ CH ₂ (CH ₃)CH-(CF ₃)CH | H |
| 85 | CH ₃ (CH ₃)CHCH ₂ (CH ₃)-CH | H |
| 86 | CH ₃ (CH ₃)CHCH ₂ (CF ₃)-CH | H |
| 87 | (S)-CH ₃ (CH ₃)CHCH ₂ -(CF ₃)CH | H |
| 88 | (R)-CH ₃ (CH ₃)CHCH ₂ -(CF ₃)CH | H |
| 89 | (CH ₃) ₂ CH(CH ₃)CH-(CH ₃)CH ₂ | H |
| 90 | (CH ₃) ₃ CCH ₂ (CH ₃)CH | H |
| 91 | <i>E</i> -CH ₃ CH=CH(CH ₃)CH | H |
| 92 | <i>E</i> -CH ₃ CH=CH(CH ₃ CH ₂)-CH | H |
| 93 | CH ₃ CH ₂ CH ₂ (CH ₃ CH ₂)CH | H |
| 94 | CH ₃ CH ₂ (CH ₃ CH ₂)CHCH ₂ | H |
| 95 | CF ₂ =CFCH ₂ CH ₂ | H |
| 96 | CF ₃ CH ₂ (CH ₃)CHCH ₂ | H |
| 97 | CF ₃ CF ₂ CH ₂ CH ₂ | H |
| 98 | CF ₃ CF ₂ CF ₂ CH ₂ | H |
| 99 | CF ₂ =C(CH ₃)CH ₂ CH ₂ | H |
| 100 | CH ₃ CH ₂ CH ₂ CH ₂ CH ₂ | H |
| 101 | CH ₃ CH ₂ CH ₂ CH ₂ (CH ₃)CH | H |
| 102 | CH ₃ CH ₂ CH ₂ (CH ₃)CHCH ₂ | H |
| 103 | CH ₃ CH ₂ (CH ₃)CHCH ₂ CH ₂ | H |

| Cmpd No | R ³ | R ⁴ |
|---------|-------------------------------------------------------------------------------------------|----------------|
| 104 | CH ₃ CH ₂ CH ₂ (CH ₃)CH-(CH ₃)CH | H |
| 105 | CH ₃ CH ₂ (CH ₃)CHCH ₂ -(CH ₃)CH | H |
| 106 | HOCH ₂ CH ₂ | H |
| 107 | CH ₃ OCH ₂ CH ₂ | H |
| 108 | CH ₃ OCH ₂ (CH ₃)CH | H |
| 109 | CH ₃ OCH ₂ (CF ₃)CH | H |
| 110 | CH ₃ OCH ₂ (CH ₃) ₂ C | H |
| 111 | CH ₃ O(CH ₃)CHCH ₂ | H |
| 112 | CH ₃ O(CH ₃)CH(CH ₃)CH | H |
| 113 | HC≡CCH ₂ | H |
| 114 | CH ₃ C≡CCH ₂ | H |
| 115 | HC≡CCH ₂ CH ₂ | H |
| 116 | HOCH ₂ CH ₂ CH ₂ | H |
| 117 | CH ₃ OCH ₂ CH ₂ CH ₂ | H |
| 118 | (CH ₃) ₃ SiCH ₂ | H |
| 119 | C ₆ H ₅ CH ₂ | H |
| 120 | C ₆ H ₅ (CH ₃)CH | H |
| 121 | 4-F-C ₆ H ₄ CH ₂ | H |
| 122 | 4-Cl-C ₆ H ₄ CH ₂ | H |
| 123 | 4-F-C ₆ H ₅ (CH ₃)CH | H |
| 124 | 4-Cl-C ₆ H ₅ (CH ₃)CH | H |
| 125 | C ₆ H ₅ CH ₂ CH ₂ | H |
| 126 | 4-F-C ₆ H ₅ CH ₂ CH ₂ | H |
| 127 | 1-piperidino | H |
| 128 | 1-pyrrolidino | H |
| 129 | cyclo-C ₅ H ₉ CH ₂ | H |
| 130 | Bicyclo[2.2.1]hept-2-yl | H |
| 131 | 1-CH ₃ -cyclopropyl | H |
| 132 | <i>cis</i> -2-CH ₃ -cyclopropyl | H |
| 133 | <i>trans</i> -2-CH ₃ -cyclopropyl | H |
| 134 | 2,2-(CH ₃) ₂ -cyclopropyl | H |
| 135 | 1-CH ₃ -cyclobutyl | H |
| 136 | <i>cis</i> -2-CH ₃ -cyclobutyl | H |
| 137 | <i>trans</i> -2-CH ₃ -cyclobutyl | H |
| 138 | <i>cis</i> -3-CH ₃ -cyclobutyl | H |

| Cmpd No | R ³ | R ⁴ |
|---------|-----------------------------------------------------------------------------------------------------|----------------|
| 139 | <i>trans</i> -3-CH ₃ -cyclobutyl | H |
| 140 | 2,2-(CH ₃) ₂ -cyclobutyl | H |
| 141 | 3,3-(CH ₃) ₂ -cyclobutyl | H |
| 142 | 1-CH ₃ -cyclopentyl | H |
| 143 | <i>cis</i> -2-CH ₃ -cyclopentyl | H |
| 144 | <i>trans</i> -2-CH ₃ -cyclopentyl | H |
| 145 | <i>cis</i> -3-CH ₃ -cyclopentyl | H |
| 146 | <i>trans</i> -3-CH ₃ -cyclopentyl | H |
| 147 | 2,2-(CH ₃) ₂ -cyclopentyl | H |
| 148 | 3,3-(CH ₃) ₂ -cyclopentyl | H |
| 149 | 1-CH ₃ -cyclohexyl | H |
| 150 | <i>cis</i> -2-CH ₃ -cyclohexyl | H |
| 151 | <i>trans</i> -2-CH ₃ -cyclohexyl | H |
| 152 | <i>cis</i> -3-CH ₃ -cyclohexyl | H |
| 153 | <i>trans</i> -3-CH ₃ -cyclohexyl | H |
| 154 | 2,2-(CH ₃) ₂ -cyclohexyl | H |
| 155 | 3,3-(CH ₃) ₂ -cyclohexyl | H |
| 156 | <i>cis</i> -4-CH ₃ -cyclohexyl | H |
| 157 | <i>trans</i> -4-CH ₃ -cyclohexyl | H |
| 158 | 4,4-(CH ₃) ₂ -cyclohexyl | H |
| 159 | 4-(CH ₃) ₃ C-cyclohexyl | H |
| 160 | -(CH ₂) ₃ - | |
| 161 | -(CH ₂) ₄ - | |
| 162 | -(CH ₂) ₅ - | |
| 163 | -(CH ₂) ₆ - | |
| 164 | -(CH ₂) ₂ (CH ₃) ₂ C(CH ₂) ₂ - | |
| 165 | -(CH ₃)CH(CH ₂) ₂ - | |
| 166 | -(CH ₃)CH(CH ₂) ₃ - | |
| 167 | -(CH ₃)CH(CH ₂) ₄ - | |
| 168 | -(CH ₃)CH(CH ₂) ₅ - | |
| 169 | -CH ₂ CH=CH(CH ₂) ₂ - | |
| 170 | -(CH ₂) ₂ NH(CH ₂) ₂ - | |
| 171 | -(CH ₂) ₂ NCH ₃ (CH ₂) ₂ - | |
| 172 | -(CH ₂) ₂ S(CH ₂) ₂ - | |
| 173 | -(CH ₂) ₂ SO(CH ₂) ₂ - | |
| 174 | -(CH ₂) ₂ SO ₂ (CH ₂) ₂ - | |

| Cmpd No | R ³ | R ⁴ |
|---------|------------------------------------------------------------------------------|-----------------|
| 175 | -CH ₂ (CH ₃)CHO(CH ₃)CHCH ₂ - | |
| 176 | C ₂ H ₅ | CH ₃ |
| 177 | <i>n</i> -C ₃ H ₇ | CH ₃ |
| 178 | <i>i</i> -C ₃ H ₇ | CH ₃ |
| 179 | <i>n</i> -C ₄ H ₉ | CH ₃ |
| 180 | <i>t</i> -C ₄ H ₉ | CH ₃ |
| 181 | CH ₂ =CHCH ₂ | CH ₃ |
| 182 | CH ₂ =C(CH ₃)CH ₂ | CH ₃ |
| 183 | CF ₃ CH ₂ | CH ₃ |
| 184 | CF ₃ CH ₂ CH ₂ | CH ₃ |
| 185 | CF ₃ CH ₂ CH ₂ CH ₂ | CH ₃ |
| 186 | CF ₃ (CH ₃)CH | CH ₃ |
| 187 | (S)-CF ₃ (CH ₃)CH | CH ₃ |
| 188 | (R)-CF ₃ (CH ₃)CH | CH ₃ |
| 189 | cyclo-C ₃ H ₅ | CH ₃ |
| 190 | cyclo-C ₄ H ₇ | CH ₃ |
| 191 | cyclo-C ₅ H ₉ | CH ₃ |
| 192 | cyclo-C ₆ H ₁₁ | CH ₃ |
| 193 | cyclo-C ₃ H ₅ CH ₂ | CH ₃ |
| 194 | cyclo-C ₄ H ₇ CH ₂ | CH ₃ |
| 195 | cyclo-C ₆ H ₁₁ CH ₂ | CH ₃ |
| 196 | CH ₃ CH ₂ (CH ₃)CH | CH ₃ |
| 197 | (S)-CH ₃ CH ₂ (CH ₃)CH | CH ₃ |
| 198 | (R)-CH ₃ CH ₂ (CH ₃)CH | CH ₃ |
| 199 | cyclo-C ₇ H ₁₃ | CH ₃ |
| 200 | CH ₂ =C(CH ₃)CH ₂ | CH ₃ |
| 201 | CF ₃ CH ₂ | CH ₃ |
| 202 | 4- <i>t</i> -C ₄ H ₉ -C ₆ H ₄ NH | CH ₃ |
| 203 | 4-F-C ₆ H ₄ NH | CH ₃ |
| 204 | C ₆ H ₅ NH | CH ₃ |
| 205 | 4-CH ₃ -C ₆ H ₄ NH | CH ₃ |
| 206 | 4-Br-C ₆ H ₄ NH | CH ₃ |
| 207 | 2-F-C ₆ H ₄ NH | CH ₃ |
| 208 | 3,4-Cl ₂ -C ₆ H ₃ NH | CH ₃ |
| 209 | 3-CF ₃ -C ₆ H ₄ NH | CH ₃ |
| 210 | 3,5-Cl ₂ -C ₆ H ₃ NH | CH ₃ |

| Cmpd No | R ³ | R ⁴ |
|---------|-----------------------------------------------------------------------------------|-----------------|
| 211 | 4-CF ₃ O-C ₆ H ₅ NH | CH ₃ |
| 212 | 2-CF ₃ -C ₆ H ₄ NH | CH ₃ |
| 213 | 4-CF ₃ -C ₆ H ₄ NH | CH ₃ |
| 214 | 2-Br-C ₆ H ₄ NH | CH ₃ |
| 215 | 2-Cl-C ₆ H ₄ NH | CH ₃ |
| 216 | 2-CH ₃ -4-Cl-C ₆ H ₃ NH | CH ₃ |
| 217 | 2-CH ₃ -5-F-C ₆ H ₃ NH | CH ₃ |
| 218 | 3-Cl-C ₆ H ₄ NH | CH ₃ |
| 219 | CH ₃ | CH ₃ |
| 220 | (CH ₃) ₂ CHCH ₂ | CH ₃ |
| 221 | (CH ₃) ₃ CCH ₂ | CH ₃ |
| 222 | (CH ₃) ₃ C(CH ₃)CH | CH ₃ |
| 223 | CH ₃ CH ₂ (CH ₃) ₂ C | CH ₃ |
| 224 | CH ₃ CH ₂ (CF ₃)CH | CH ₃ |
| 225 | (S)-CH ₃ CH ₂ (CF ₃)CH | CH ₃ |
| 226 | (R)-CH ₃ CH ₂ (CF ₃)CH | CH ₃ |
| 227 | CH ₃ CH ₂ (CH ₃ CH ₂)CH | CH ₃ |
| 228 | (CH ₃) ₂ CH(CH ₃ CH ₂)CH | CH ₃ |
| 229 | (CH ₃) ₂ CH(CH ₃)CH | CH ₃ |
| 230 | (CH ₃) ₂ CH(CF ₃)CH | CH ₃ |
| 231 | (S)-(CH ₃) ₂ CH(CF ₃)CH | CH ₃ |
| 232 | (R)-(CH ₃) ₂ CH(CF ₃)CH | CH ₃ |
| 233 | HC=C(CH ₃)CH ₂ | CH ₃ |
| 234 | CH ₂ =CH(CH ₃ CH ₂)CH | CH ₃ |
| 235 | CH ₃ CH ₂ CH ₂ (CH ₃)CH | CH ₃ |
| 236 | CH ₃ CH ₂ CH ₂ (CF ₃)CH | CH ₃ |
| 237 | (S)-CH ₃ CH ₂ CH ₂ (CF ₃)CH | CH ₃ |
| 238 | (R)-CH ₃ CH ₂ CH ₂ (CF ₃)CH | CH ₃ |
| 239 | CH ₃ CH ₂ CH ₂ (CH ₃) ₂ C | CH ₃ |
| 240 | CH ₃ CH ₂ (CH ₃)CHCH ₂ | CH ₃ |
| 241 | (CH ₃) ₂ CHCH ₂ CH ₂ | CH ₃ |
| 242 | (CH ₃) ₃ CCH ₂ CH ₂ | CH ₃ |
| 243 | CH ₃ CH ₂ (CH ₃)CH(CH ₃)-CH | CH ₃ |
| 244 | CH ₃ CH ₂ (CH ₃)CH(CF ₃)-CH | CH ₃ |
| 245 | (S)-CH ₃ CH ₂ (CH ₃)CH-(CF ₃)CH | CH ₃ |

| Cmpd No | R ³ | R ⁴ |
|---------|-------------------------------------------------------------------------------------------|-----------------|
| 246 | (R)-CH ₃ CH ₂ (CH ₃)CH-(CF ₃)CH | CH ₃ |
| 247 | CH ₃ (CH ₃)CHCH ₂ -(CH ₃)CH | CH ₃ |
| 248 | CH ₃ (CH ₃)CHCH ₂ -(CF ₃)CH | CH ₃ |
| 249 | (S)-CH ₃ (CH ₃)CHCH ₂ -(CF ₃)CH | CH ₃ |
| 250 | (R)-CH ₃ (CH ₃)HCH ₂ -(CF ₃)CH | CH ₃ |
| 251 | (CH ₃) ₂ CH(CH ₃)CH(CH ₃)-CH ₂ | CH ₃ |
| 252 | (CH ₃) ₃ CCH ₂ (CH ₃)CH | CH ₃ |
| 253 | E-CH ₃ CH=CH(CH ₃)CH | CH ₃ |
| 254 | E-CH ₃ CH=CH(CH ₃ CH ₂)-CH | CH ₃ |
| 255 | CH ₃ CH ₂ CH ₂ (CH ₃ CH ₂)-CH | CH ₃ |
| 256 | CH ₃ CH ₂ (CH ₃ CH ₂)CHCH ₂ | CH ₃ |
| 257 | CF ₂ =CFCH ₂ CH ₂ | CH ₃ |
| 258 | CF ₃ CH ₂ (CH ₃)CHCH ₂ | CH ₃ |
| 259 | CF ₃ CF ₂ CH ₂ CH ₂ | CH ₃ |
| 260 | CF ₃ CF ₂ CF ₂ CH ₂ | CH ₃ |
| 261 | CF ₂ =C(CH ₃)CH ₂ CH ₂ | CH ₃ |
| 262 | CH ₃ CH ₂ CH ₂ CH ₂ CH ₂ | CH ₃ |
| 263 | CH ₃ CH ₂ CH ₂ CH ₂ (CH ₃)CH | CH ₃ |
| 264 | CH ₃ CH ₂ CH ₂ (CH ₃)CHCH ₂ | CH ₃ |
| 265 | CH ₃ CH ₂ (CH ₃)CHCH ₂ CH ₂ | CH ₃ |
| 266 | CH ₃ CH ₂ CH ₂ (CH ₃)CH-(CH ₃)CH | CH ₃ |
| 267 | CH ₃ CH ₂ (CH ₃)CHCH ₂ -(CH ₃)CH | CH ₃ |
| 268 | HOCH ₂ CH ₂ | CH ₃ |
| 269 | CH ₃ OCH ₂ CH ₂ | CH ₃ |
| 270 | CH ₃ OCH ₂ (CH ₃)CH | CH ₃ |
| 271 | CH ₃ OCH ₂ (CF ₃)CH | CH ₃ |
| 272 | CH ₃ OCH ₂ (CH ₃) ₂ C | CH ₃ |
| 273 | CH ₃ O(CH ₃)CHCH ₂ | CH ₃ |
| 274 | CH ₃ O(CH ₃)CH(CH ₃)CH | CH ₃ |
| 275 | HC≡CCH ₂ | CH ₃ |
| 276 | CH ₃ C≡CCH ₂ | CH ₃ |
| 277 | HC≡CCH ₂ CH ₂ | CH ₃ |
| 278 | HOCH ₂ CH ₂ CH ₂ | CH ₃ |

| Cmpd No | R ³ | R ⁴ |
|---------|-------------------------------------------------------------------|-----------------|
| 279 | CH ₃ OCH ₂ CH ₂ CH ₂ | CH ₃ |
| 280 | (CH ₃) ₃ SiCH ₂ | CH ₃ |
| 281 | C ₆ H ₅ CH ₂ | CH ₃ |
| 282 | C ₆ H ₅ (CH ₃)CH | CH ₃ |
| 283 | 4-F-C ₆ H ₄ CH ₂ | H |
| 284 | 4-Cl-C ₆ H ₄ CH ₂ | CH ₃ |
| 285 | 4-F-C ₆ H ₅ (CH ₃)CH | CH ₃ |
| 286 | 4-Cl-C ₆ H ₅ (CH ₃)CH | CH ₃ |
| 287 | C ₆ H ₅ CH ₂ CH ₂ | CH ₃ |
| 288 | 4-F-C ₆ H ₅ CH ₂ CH ₂ | CH ₃ |
| 289 | 1-piperidino | CH ₃ |
| 290 | 1-pyrrolidino | CH ₃ |
| 291 | cyclo-C ₃ H ₅ CH ₂ | CH ₃ |
| 292 | bicyclo[2.2.1]hept-2-yl | CH ₃ |
| 293 | 1-CH ₃ -cyclopropyl | CH ₃ |
| 294 | <i>cis</i> -2-CH ₃ -cyclopropyl | CH ₃ |
| 295 | <i>trans</i> -2-CH ₃ -cyclopropyl | CH ₃ |
| 296 | 2,2-(CH ₃) ₂ -cyclopropyl | CH ₃ |
| 297 | 1-CH ₃ -cyclobutyl | CH ₃ |
| 298 | <i>cis</i> -2-CH ₃ -cyclobutyl | CH ₃ |
| 299 | <i>trans</i> -2-CH ₃ -cyclobutyl | CH ₃ |
| 300 | <i>cis</i> -3-CH ₃ -cyclobutyl | CH ₃ |
| 301 | <i>trans</i> -3-CH ₃ -cyclobutyl | CH ₃ |
| 302 | 2,2-(CH ₃) ₂ -cyclobutyl | CH ₃ |
| 303 | 3,3-(CH ₃) ₂ -cyclobutyl | CH ₃ |
| 304 | 1-CH ₃ -cyclopentyl | CH ₃ |
| 305 | <i>cis</i> -2-CH ₃ -cyclopentyl | CH ₃ |
| 306 | <i>trans</i> -2-CH ₃ -cyclopentyl | CH ₃ |
| 307 | <i>cis</i> -3-CH ₃ -cyclopentyl | CH ₃ |
| 308 | <i>trans</i> -3-CH ₃ -cyclopentyl | CH ₃ |
| 309 | 2,2-(CH ₃) ₂ -cyclopentyl | CH ₃ |
| 310 | 3,3-(CH ₃) ₂ -cyclopentyl | CH ₃ |
| 311 | 1-CH ₃ -cyclohexyl | CH ₃ |
| 312 | <i>cis</i> -2-CH ₃ -cyclohexyl | CH ₃ |
| 313 | <i>trans</i> -2-CH ₃ -cyclohexyl | CH ₃ |
| 314 | <i>cis</i> -3-CH ₃ -cyclohexyl | CH ₃ |

| Cmpd No | R ³ | R ⁴ |
|---------|------------------------------------------------------------------------------|-------------------------------|
| 315 | <i>trans</i> -3-CH ₃ -cyclohexyl | CH ₃ |
| 316 | 2,2-(CH ₃) ₂ -cyclohexyl | CH ₃ |
| 317 | 3,3-(CH ₃) ₂ -cyclohexyl | CH ₃ |
| 318 | <i>cis</i> -4-CH ₃ -cyclohexyl | CH ₃ |
| 319 | <i>trans</i> -4-CH ₃ -cyclohexyl | CH ₃ |
| 320 | 4,4-(CH ₃) ₂ -cyclohexyl | CH ₃ |
| 321 | 4-(CH ₃) ₃ C-cyclohexyl | CH ₃ |
| 322 | <i>n</i> -C ₃ H ₇ | C ₂ H ₅ |
| 323 | <i>i</i> -C ₃ H ₇ | C ₂ H ₅ |
| 324 | <i>n</i> -C ₄ H ₉ | C ₂ H ₅ |
| 325 | <i>t</i> -C ₄ H ₉ | C ₂ H ₅ |
| 326 | CH ₂ =CHCH ₂ | C ₂ H ₅ |
| 327 | CF ₃ CH ₂ CH ₂ | C ₂ H ₅ |
| 328 | CF ₃ CH ₂ CH ₂ CH ₂ | C ₂ H ₅ |
| 329 | CF ₃ (CH ₃)CH | C ₂ H ₅ |
| 330 | (S)-CF ₃ (CH ₃)CH | C ₂ H ₅ |
| 331 | (R)-CF ₃ (CH ₃)CH | C ₂ H ₅ |
| 332 | cyclo-C ₃ H ₅ | C ₂ H ₅ |
| 333 | cyclo-C ₄ H ₇ | C ₂ H ₅ |
| 334 | cyclo-C ₅ H ₉ | C ₂ H ₅ |
| 335 | cyclo-C ₆ H ₁₁ | C ₂ H ₅ |
| 336 | cyclo-C ₃ H ₅ CH ₂ | C ₂ H ₅ |
| 337 | cyclo-C ₄ H ₇ CH ₂ | C ₂ H ₅ |
| 338 | cyclo-C ₆ H ₁₁ CH ₂ | C ₂ H ₅ |
| 339 | CH ₃ CH ₂ (CH ₃)CH | C ₂ H ₅ |
| 340 | (S)-CH ₃ CH ₂ (CH ₃)CH | C ₂ H ₅ |
| 341 | (R)-CH ₃ CH ₂ (CH ₃)CH | C ₂ H ₅ |
| 342 | cyclo-C ₇ H ₁₃ | C ₂ H ₅ |
| 343 | 4- <i>t</i> -C ₄ H ₉ -C ₆ H ₄ NH | C ₂ H ₅ |
| 344 | 4-F-C ₆ H ₄ NH | C ₂ H ₅ |
| 345 | C ₆ H ₅ NH | C ₂ H ₅ |
| 346 | 4-CH ₃ -C ₆ H ₄ NH | C ₂ H ₅ |
| 347 | 4-Br-C ₆ H ₄ NH | C ₂ H ₅ |
| 348 | 2-F-C ₆ H ₄ NH | C ₂ H ₅ |
| 349 | 3,4-Cl ₂ -C ₆ H ₃ NH | C ₂ H ₅ |
| 350 | 3-CF ₃ -C ₆ H ₄ NH | C ₂ H ₅ |

| Cmpd No | R ³ | R ⁴ |
|---------|-----------------------------------------------------------------------------------|-------------------------------|
| 351 | 3,5-Cl ₂ -C ₆ H ₃ NH | C ₂ H ₅ |
| 352 | 4-CF ₃ O-C ₆ H ₅ NH | C ₂ H ₅ |
| 353 | 2-CF ₃ -C ₆ H ₄ NH | C ₂ H ₅ |
| 354 | 4-CF ₃ -C ₆ H ₄ NH | C ₂ H ₅ |
| 355 | 2-Br-C ₆ H ₄ NH | C ₂ H ₅ |
| 356 | 2-Cl-C ₆ H ₄ NH | C ₂ H ₅ |
| 357 | 2-CH ₃ -4-Cl-C ₆ H ₃ NH | C ₂ H ₅ |
| 358 | 2-CH ₃ -5-F-C ₆ H ₃ NH | C ₂ H ₅ |
| 359 | 3-Cl-C ₆ H ₄ NH | C ₂ H ₅ |
| 360 | (CH ₃) ₂ CHCH ₂ | C ₂ H ₅ |
| 361 | (CH ₃) ₃ CCH ₂ | C ₂ H ₅ |
| 362 | (CH ₃) ₃ C(CH ₃)CH | C ₂ H ₅ |
| 363 | CH ₃ CH ₂ (CH ₃) ₂ C | C ₂ H ₅ |
| 364 | CH ₃ CH ₂ (CF ₃)CH | C ₂ H ₅ |
| 365 | (S)-CH ₃ CH ₂ (CF ₃)CH | C ₂ H ₅ |
| 366 | (R)-CH ₃ CH ₂ (CF ₃)CH | C ₂ H ₅ |
| 367 | CH ₃ CH ₂ (CH ₃ CH ₂)CH | C ₂ H ₅ |
| 368 | (CH ₃) ₂ CH(CH ₃ CH ₂)CH | C ₂ H ₅ |
| 369 | (CH ₃) ₂ CH(CH ₃)CH | C ₂ H ₅ |
| 370 | (CH ₃) ₂ CH(CF ₃)CH | C ₂ H ₅ |
| 371 | (S)-(CH ₃) ₂ CH(CF ₃)CH | C ₂ H ₅ |
| 372 | (R)-(CH ₃) ₂ CH(CF ₃)CH | C ₂ H ₅ |
| 373 | HC=C(CH ₃)CH ₂ | C ₂ H ₅ |
| 374 | CH ₂ =CH(CH ₃ CH ₂)CH | C ₂ H ₅ |
| 375 | CH ₃ CH ₂ CH ₂ (CH ₃)CH | C ₂ H ₅ |
| 376 | CH ₃ CH ₂ CH ₂ (CF ₃)CH | C ₂ H ₅ |
| 377 | (S)-CH ₃ CH ₂ CH ₂ (CF ₃)CH | C ₂ H ₅ |
| 378 | (R)-CH ₃ CH ₂ CH ₂ (CF ₃)CH | C ₂ H ₅ |
| 379 | CH ₃ CH ₂ CH ₂ (CH ₃) ₂ C | C ₂ H ₅ |
| 380 | CH ₃ CH ₂ (CH ₃)CHCH ₂ | C ₂ H ₅ |
| 381 | (CH ₃) ₂ CHCH ₂ CH ₂ | C ₂ H ₅ |
| 382 | (CH ₃) ₃ CCH ₂ CH ₂ | C ₂ H ₅ |
| 383 | CH ₃ CH ₂ (CH ₃)CH(CH ₃)-CH | C ₂ H ₅ |
| 384 | CH ₃ CH ₂ (CH ₃)CH(CF ₃)-CH | C ₂ H ₅ |
| 385 | (S)-CH ₃ CH ₂ (CH ₃)CH-(CF ₃)CH | C ₂ H ₅ |

| Cmpd No | R ³ | R ⁴ |
|---------|-------------------------------------------------------------------------------------------|-------------------------------|
| 386 | (R)-CH ₃ CH ₂ (CH ₃)CH-(CF ₃)CH | C ₂ H ₅ |
| 387 | CH ₃ (CH ₃)CHCH ₂ (CH ₃)-CH | C ₂ H ₅ |
| 388 | CH ₃ (CH ₃)CHCH ₂ (CF ₃)-CH | C ₂ H ₅ |
| 389 | (S)-CH ₃ (CH ₃)CHCH ₂ -(CF ₃)CH | C ₂ H ₅ |
| 390 | (R)-CH ₃ (CH ₃)CHCH ₂ -(CF ₃)CH | C ₂ H ₅ |
| 391 | (CH ₃) ₂ CH(CH ₃)CH-(CH ₃)CH ₂ | C ₂ H ₅ |
| 392 | (CH ₃) ₃ CCH ₂ (CH ₃)CH | C ₂ H ₅ |
| 393 | E-CH ₃ CH=CH(CH ₃)CH | C ₂ H ₅ |
| 394 | E-C H ₃ CH=CH-(CH ₃ CH ₂)CH | C ₂ H ₅ |
| 395 | CH ₃ CH ₂ CH ₂ (CH ₃ CH ₂)CH | C ₂ H ₅ |
| 396 | CH ₃ CH ₂ (CH ₃ CH ₂)CHCH ₂ | C ₂ H ₅ |
| 397 | CF ₂ =CFCH ₂ CH ₂ | C ₂ H ₅ |
| 398 | CF ₃ CH ₂ (CH ₃)CHCH ₂ | C ₂ H ₅ |
| 399 | CF ₃ CF ₂ CH ₂ CH ₂ | C ₂ H ₅ |
| 400 | CF ₃ CF ₂ CF ₂ CH ₂ | C ₂ H ₅ |
| 401 | CF ₂ =C(CH ₃)CH ₂ CH ₂ | C ₂ H ₅ |
| 402 | CH ₃ CH ₂ CH ₂ CH ₂ CH ₂ | C ₂ H ₅ |
| 403 | CH ₃ CH ₂ CH ₂ CH ₂ (CH ₃)CH | C ₂ H ₅ |
| 404 | CH ₃ CH ₂ CH ₂ (CH ₃)CHCH ₂ | C ₂ H ₅ |
| 405 | CH ₃ CH ₂ (CH ₃)CHCH ₂ CH ₂ | C ₂ H ₅ |
| 406 | CH ₃ CH ₂ CH ₂ (CH ₃)CH-(CH ₃)CH | C ₂ H ₅ |
| 407 | CH ₃ CH ₂ (CH ₃)CHCH ₂ -(CH ₃)CH | C ₂ H ₅ |
| 408 | HOCH ₂ CH ₂ | C ₂ H ₅ |
| 409 | CH ₃ OCH ₂ CH ₂ | C ₂ H ₅ |
| 410 | CH ₃ OCH ₂ (CH ₃)CH | C ₂ H ₅ |
| 411 | CH ₃ OCH ₂ (CF ₃)CH | C ₂ H ₅ |
| 412 | CH ₃ OCH ₂ (CH ₃) ₂ C | C ₂ H ₅ |
| 413 | CH ₃ O(CH ₃)CHCH ₂ | C ₂ H ₅ |
| 414 | CH ₃ O(CH ₃)CH(CH ₃)CH | C ₂ H ₅ |
| 415 | HC≡CCH ₂ | C ₂ H ₅ |
| 416 | CH ₃ C≡CCH ₂ | C ₂ H ₅ |
| 417 | HC≡CCH ₂ CH ₂ | C ₂ H ₅ |
| 418 | HOCH ₂ CH ₂ CH ₂ | C ₂ H ₅ |

| Cmpd No | R ³ | R ⁴ |
|---------|-------------------------------------------------------------------|-------------------------------|
| 419 | CH ₃ OCH ₂ CH ₂ CH ₂ | C ₂ H ₅ |
| 420 | (CH ₃) ₃ SiCH ₂ | C ₂ H ₅ |
| 421 | C ₆ H ₅ CH ₂ | C ₂ H ₅ |
| 422 | C ₆ H ₅ (CH ₃)CH | C ₂ H ₅ |
| 423 | 4-F-C ₆ H ₄ CH ₂ | C ₂ H ₅ |
| 424 | 4-Cl-C ₆ H ₄ CH ₂ | C ₂ H ₅ |
| 425 | 4-F-C ₆ H ₅ (CH ₃)CH | C ₂ H ₅ |
| 426 | 4-Cl-C ₆ H ₅ (CH ₃)CH | C ₂ H ₅ |
| 427 | C ₆ H ₅ CH ₂ CH ₂ | C ₂ H ₅ |
| 428 | 4-F-C ₆ H ₅ CH ₂ CH ₂ | C ₂ H ₅ |
| 429 | 1-piperidino | C ₂ H ₅ |
| 430 | 1-pyrrolidino | C ₂ H ₅ |
| 431 | cyclo-C ₃ H ₅ CH ₂ | C ₂ H ₅ |
| 432 | Bicyclo[2.2.1]hept-2-yl | C ₂ H ₅ |
| 433 | 1-CH ₃ -cyclopropyl | C ₂ H ₅ |
| 434 | <i>cis</i> -2-CH ₃ -cyclopropyl | C ₂ H ₅ |
| 435 | <i>trans</i> -2-CH ₃ -cyclopropyl | C ₂ H ₅ |
| 436 | 2,2-(CH ₃) ₂ -cyclopropyl | C ₂ H ₅ |
| 437 | 1-CH ₃ -cyclobutyl | C ₂ H ₅ |
| 438 | <i>cis</i> -2-CH ₃ -cyclobutyl | C ₂ H ₅ |
| 439 | <i>trans</i> -2-CH ₃ -cyclobutyl | C ₂ H ₅ |
| 440 | <i>cis</i> -3-CH ₃ -cyclobutyl | C ₂ H ₅ |
| 441 | <i>trans</i> -3-CH ₃ -cyclobutyl | C ₂ H ₅ |
| 442 | 2,2-(CH ₃) ₂ -cyclobutyl | C ₂ H ₅ |
| 443 | 3,3-(CH ₃) ₂ -cyclobutyl | C ₂ H ₅ |
| 444 | 1-CH ₃ -cyclopentyl | C ₂ H ₅ |
| 445 | <i>cis</i> -2-CH ₃ -cyclopentyl | C ₂ H ₅ |
| 446 | <i>trans</i> -2-CH ₃ -cyclopentyl | C ₂ H ₅ |
| 447 | <i>cis</i> -3-CH ₃ -cyclopentyl | C ₂ H ₅ |
| 448 | <i>trans</i> -3-CH ₃ -cyclopentyl | C ₂ H ₅ |
| 449 | 2,2-(CH ₃) ₂ -cyclopentyl | C ₂ H ₅ |
| 450 | 3,3-(CH ₃) ₂ -cyclopentyl | C ₂ H ₅ |
| 451 | 1-CH ₃ -cyclohexyl | C ₂ H ₅ |
| 452 | <i>cis</i> -2-CH ₃ -cyclohexyl | C ₂ H ₅ |
| 453 | <i>trans</i> -2-CH ₃ -cyclohexyl | C ₂ H ₅ |
| 454 | <i>cis</i> -3-CH ₃ -cyclohexyl | C ₂ H ₅ |

| Cmpd No | R ³ | R ⁴ |
|---------|------------------------------------------------------------------------------|---------------------------------|
| 355 | <i>trans</i> -3-CH ₃ -cyclohexyl | C ₂ H ₅ |
| 456 | 2,2-(CH ₃) ₂ -cyclohexyl | C ₂ H ₅ |
| 457 | 3,3-(CH ₃) ₂ -cyclohexyl | C ₂ H ₅ |
| 458 | <i>cis</i> -4-CH ₃ -cyclohexyl | C ₂ H ₅ |
| 459 | <i>trans</i> -4-CH ₃ -cyclohexyl | C ₂ H ₅ |
| 460 | 4,4-(CH ₃) ₂ -cyclohexyl | C ₂ H ₅ |
| 461 | 4-(CH ₃) ₃ C-cyclohexyl | C ₂ H ₅ |
| 462 | <i>n</i> -C ₃ H ₇ | CF ₃ CH ₂ |
| 463 | <i>i</i> -C ₃ H ₇ | CF ₃ CH ₂ |
| 464 | <i>n</i> -C ₄ H ₉ | CF ₃ CH ₂ |
| 465 | <i>t</i> -C ₄ H ₉ | CF ₃ CH ₂ |
| 466 | CH ₂ =CHCH ₂ | CF ₃ CH ₂ |
| 467 | CH ₂ =C(CH ₃)CH ₂ | CF ₃ CH ₂ |
| 468 | CF ₃ CH ₂ | CF ₃ CH ₂ |
| 469 | CF ₃ CH ₂ CH ₂ | CF ₃ CH ₂ |
| 470 | CF ₃ CH ₂ CH ₂ CH ₂ | CF ₃ CH ₂ |
| 471 | CF ₃ (CH ₃)CH | CF ₃ CH ₂ |
| 472 | (S)-CF ₃ (CH ₃)CH | CF ₃ CH ₂ |
| 473 | (R)-CF ₃ (CH ₃)CH | CF ₃ CH ₂ |
| 474 | cyclo-C ₃ H ₅ | CF ₃ CH ₂ |
| 475 | cyclo-C ₄ H ₇ | CF ₃ CH ₂ |
| 476 | cyclo-C ₅ H ₉ | CF ₃ CH ₂ |
| 477 | cyclo-C ₆ H ₁₁ | CF ₃ CH ₂ |
| 478 | cyclo-C ₃ H ₅ CH ₂ | CF ₃ CH ₂ |
| 479 | cyclo-C ₄ H ₇ CH ₂ | CF ₃ CH ₂ |
| 480 | cyclo-C ₆ H ₁₁ CH ₂ | CF ₃ CH ₂ |
| 481 | CH ₃ CH ₂ (CH ₃)CH | CF ₃ CH ₂ |
| 482 | (S)-CH ₃ CH ₂ (CH ₃)CH | CF ₃ CH ₂ |
| 483 | (R)-CH ₃ CH ₂ (CH ₃)CH | CF ₃ CH ₂ |
| 484 | cyclo-C ₇ H ₁₃ | CF ₃ CH ₂ |
| 485 | CH ₂ =C(CH ₃)CH ₂ | CF ₃ CH ₂ |
| 486 | CF ₃ CH ₂ | CF ₃ CH ₂ |
| 487 | 4- <i>t</i> -C ₄ H ₉ -C ₆ H ₄ NH | CF ₃ CH ₂ |
| 488 | 4-F-C ₆ H ₄ NH | CF ₃ CH ₂ |
| 489 | C ₆ H ₅ NH | CF ₃ CH ₂ |
| 490 | 4-CH ₃ -C ₆ H ₄ NH | CF ₃ CH ₂ |

| Cmpd No | R ³ | R ⁴ |
|---------|-----------------------------------------------------------------------------------|---------------------------------|
| 491 | 4-Br-C ₆ H ₄ NH | CF ₃ CH ₂ |
| 492 | 2-F-C ₆ H ₄ NH | CF ₃ CH ₂ |
| 493 | 3,4-Cl ₂ -C ₆ H ₃ NH | CF ₃ CH ₂ |
| 494 | 3-CF ₃ -C ₆ H ₄ NH | CF ₃ CH ₂ |
| 495 | 3,5-Cl ₂ -C ₆ H ₃ NH | CF ₃ CH ₂ |
| 496 | 4-CF ₃ O-C ₆ H ₃ NH | CF ₃ CH ₂ |
| 497 | 2-CF ₃ -C ₆ H ₄ NH | CF ₃ CH ₂ |
| 498 | 4-CF ₃ -C ₆ H ₄ NH | CF ₃ CH ₂ |
| 499 | 2-Br-C ₆ H ₄ NH | CF ₃ CH ₂ |
| 500 | 2-Cl-C ₆ H ₄ NH | CF ₃ CH ₂ |
| 501 | 2-CH ₃ -4-Cl-C ₆ H ₃ NH | CF ₃ CH ₂ |
| 502 | 2-CH ₃ -5-F-C ₆ H ₃ NH | CF ₃ CH ₂ |
| 503 | 3-Cl-C ₆ H ₄ NH | CF ₃ CH ₂ |
| 504 | (CH ₃) ₂ CHCH ₂ | CF ₃ CH ₂ |
| 505 | (CH ₃) ₃ CCH ₂ | CF ₃ CH ₂ |
| 506 | (CH ₃) ₃ C(CH ₃)CH | CF ₃ CH ₂ |
| 507 | CH ₃ CH ₂ (CH ₃) ₂ C | CF ₃ CH ₂ |
| 508 | CH ₃ CH ₂ (CF ₃)CH | CF ₃ CH ₂ |
| 509 | (S)-CH ₃ CH ₂ (CF ₃)CH | CF ₃ CH ₂ |
| 510 | (R)-CH ₃ CH ₂ (CF ₃)CH | CF ₃ CH ₂ |
| 511 | CH ₃ CH ₂ (CH ₃ CH ₂)CH | CF ₃ CH ₂ |
| 512 | (CH ₃) ₂ CH(CH ₃ CH ₂)CH | CF ₃ CH ₂ |
| 513 | (CH ₃) ₂ CH(CH ₃)CH | CF ₃ CH ₂ |
| 514 | (CH ₃) ₂ CH(CF ₃)CH | CF ₃ CH ₂ |
| 515 | (S)-(CH ₃) ₂ CH(CF ₃)CH | CF ₃ CH ₂ |
| 516 | (R)-(CH ₃) ₂ CH(CF ₃)CH | CF ₃ CH ₂ |
| 517 | HC=C(CH ₃)CH ₂ | CF ₃ CH ₂ |
| 518 | CH ₂ =CH(CH ₃ CH ₂)CH | CF ₃ CH ₂ |
| 520 | CH ₃ CH ₂ CH ₂ (CH ₃)CH | CF ₃ CH ₂ |
| 521 | CH ₃ CH ₂ CH ₂ (CF ₃)CH | CF ₃ CH ₂ |
| 522 | (S)-CH ₃ CH ₂ CH ₂ (CF ₃)CH | CF ₃ CH ₂ |
| 523 | (R)-CH ₃ CH ₂ CH ₂ (CF ₃)CH | CF ₃ CH ₂ |
| 524 | CH ₃ CH ₂ CH ₂ (CH ₃) ₂ C | CF ₃ CH ₂ |
| 525 | CH ₃ CH ₂ (CH ₃)CHCH ₂ | CF ₃ CH ₂ |
| 526 | (CH ₃) ₂ CHCH ₂ CH ₂ | CF ₃ CH ₂ |
| 527 | (CH ₃) ₃ CCH ₂ CH ₂ | CF ₃ CH ₂ |

| Cmpd No | R ³ | R ⁴ |
|---------|-------------------------------------------------------------------------------------------|---------------------------------|
| 528 | CH ₃ CH ₂ (CH ₃)CH(CH ₃)-CH | CF ₃ CH ₂ |
| 529 | CH ₃ CH ₂ (CH ₃)CH(CF ₃)-CH | CF ₃ CH ₂ |
| 530 | (S)-CH ₃ CH ₂ (CH ₃)CH-(CF ₃)CH | CF ₃ CH ₂ |
| 531 | (R)-CH ₃ CH ₂ (CH ₃)CH-(CF ₃)CH | CF ₃ CH ₂ |
| 532 | CH ₃ (CH ₃)CHCH ₂ (CH ₃)-CH | CF ₃ CH ₂ |
| 533 | CH ₃ (CH ₃)CHCH ₂ (CF ₃)-CH | CF ₃ CH ₂ |
| 534 | (S)-CH ₃ (CH ₃)CHCH ₂ -(CF ₃)CH | CF ₃ CH ₂ |
| 535 | (R)-CH ₃ (CH ₃)CHCH ₂ -(CF ₃)CH | CF ₃ CH ₂ |
| 536 | (CH ₃) ₂ CH(CH ₃)CH-(CH ₃)CH ₂ | CF ₃ CH ₂ |
| 537 | (CH ₃) ₃ CCH ₂ (CH ₃)CH | CF ₃ CH ₂ |
| 538 | <i>E</i> -CH ₃ CH=CH(CH ₃)CH | CF ₃ CH ₂ |
| 539 | <i>E</i> -CH ₃ CH=CH(CH ₃ CH ₂)-CH | CF ₃ CH ₂ |
| 540 | CH ₃ CH ₂ CH ₂ (CH ₃ CH ₂)CH | CF ₃ CH ₂ |
| 541 | CH ₃ CH ₂ (CH ₃ CH ₂)CHCH ₂ | CF ₃ CH ₂ |
| 542 | CF ₂ =CFCH ₂ CH ₂ | CF ₃ CH ₂ |
| 543 | CF ₃ CH ₂ (CH ₃)CHCH ₂ | CF ₃ CH ₂ |
| 544 | CF ₃ CF ₂ CH ₂ CH ₂ | CF ₃ CH ₂ |
| 545 | CF ₃ CF ₂ CF ₂ CH ₂ | CF ₃ CH ₂ |
| 546 | CF ₂ =C(CH ₃)CH ₂ CH ₂ | CF ₃ CH ₂ |
| 547 | CH ₃ CH ₂ CH ₂ CH ₂ CH ₂ | CF ₃ CH ₂ |
| 548 | CH ₃ CH ₂ CH ₂ CH ₂ (CH ₃)CH | CF ₃ CH ₂ |
| 549 | CH ₃ CH ₂ CH ₂ (CH ₃)CHCH ₂ | CF ₃ CH ₂ |
| 550 | CH ₃ CH ₂ (CH ₃)CHCH ₂ CH ₂ | CF ₃ CH ₂ |
| 551 | CH ₃ CH ₂ CH ₂ (CH ₃)CH-(CH ₃)CH | CF ₃ CH ₂ |
| 552 | CH ₃ CH ₂ (CH ₃)CHCH ₂ -(CH ₃)CH | CF ₃ CH ₂ |
| 553 | HOCH ₂ CH ₂ | CF ₃ CH ₂ |
| 554 | CH ₃ OCH ₂ CH ₂ | CF ₃ CH ₂ |
| 555 | CH ₃ OCH ₂ (CH ₃)CH | CF ₃ CH ₂ |
| 556 | CH ₃ OCH ₂ (CF ₃)CH | CF ₃ CH ₂ |
| 557 | CH ₃ OCH ₂ (CH ₃) ₂ C | CF ₃ CH ₂ |
| 558 | CH ₃ O(CH ₃)CHCH ₂ | CF ₃ CH ₂ |
| 559 | CH ₃ O(CH ₃)CH(CH ₃)CH | CF ₃ CH ₂ |

| Cmpd No | R ³ | R ⁴ |
|---------|-------------------------------------------------------------------|---------------------------------|
| 560 | HC≡CCH ₂ | CF ₃ CH ₂ |
| 561 | CH ₃ C≡CCH ₂ | CF ₃ CH ₂ |
| 562 | HC≡CCH ₂ CH ₂ | CF ₃ CH ₂ |
| 563 | HOCH ₂ CH ₂ CH ₂ | CF ₃ CH ₂ |
| 564 | CH ₃ OCH ₂ CH ₂ CH ₂ | CF ₃ CH ₂ |
| 565 | (CH ₃) ₃ SiCH ₂ | CF ₃ CH ₂ |
| 566 | C ₆ H ₅ CH ₂ | CF ₃ CH ₂ |
| 567 | C ₆ H ₅ (CH ₃)CH | CF ₃ CH ₂ |
| 568 | 4-F-C ₆ H ₄ CH ₂ | CF ₃ CH ₂ |
| 569 | 4-Cl-C ₆ H ₄ CH ₂ | CF ₃ CH ₂ |
| 570 | 4-F-C ₆ H ₅ (CH ₃)CH | CF ₃ CH ₂ |
| 571 | 4-Cl-C ₆ H ₅ (CH ₃)CH | CF ₃ CH ₂ |
| 572 | C ₆ H ₅ CH ₂ CH ₂ | CF ₃ CH ₂ |
| 573 | 4-F-C ₆ H ₅ CH ₂ CH ₂ | CF ₃ CH ₂ |
| 574 | 1-piperidino | CF ₃ CH ₂ |
| 575 | 1-pyrrolidino | CF ₃ CH ₂ |
| 576 | cyclo-C ₅ H ₉ CH ₂ | CF ₃ CH ₂ |
| 577 | bicyclo[2.2.1]hept-2-yl | CF ₃ CH ₂ |
| 578 | 1-CH ₃ -cyclopropyl | CF ₃ CH ₂ |
| 579 | <i>cis</i> -2-CH ₃ -cyclopropyl | CF ₃ CH ₂ |
| 580 | <i>trans</i> -2-CH ₃ -cyclopropyl | CF ₃ CH ₂ |
| 581 | 2,2-(CH ₃) ₂ -cyclopropyl | CF ₃ CH ₂ |
| 582 | 1-CH ₃ -cyclobutyl | CF ₃ CH ₂ |
| 583 | <i>cis</i> -2-CH ₃ -cyclobutyl | CF ₃ CH ₂ |
| 584 | <i>trans</i> -2-CH ₃ -cyclobutyl | CF ₃ CH ₂ |
| 585 | <i>cis</i> -3-CH ₃ -cyclobutyl | CF ₃ CH ₂ |
| 586 | <i>trans</i> -3-CH ₃ -cyclobutyl | CF ₃ CH ₂ |
| 587 | 2,2-(CH ₃) ₂ -cyclobutyl | CF ₃ CH ₂ |
| 588 | 3,3-(CH ₃) ₂ -cyclobutyl | CF ₃ CH ₂ |
| 589 | 1-CH ₃ -cyclopentyl | CF ₃ CH ₂ |
| 590 | <i>cis</i> -2-CH ₃ -cyclopentyl | CF ₃ CH ₂ |
| 591 | <i>trans</i> -2-CH ₃ -cyclopentyl | CF ₃ CH ₂ |
| 592 | <i>cis</i> -3-CH ₃ -cyclopentyl | CF ₃ CH ₂ |
| 593 | <i>trans</i> -3-CH ₃ -cyclopentyl | CF ₃ CH ₂ |
| 594 | 2,2-(CH ₃) ₂ -cyclopentyl | CF ₃ CH ₂ |
| 595 | 3,3-(CH ₃) ₂ -cyclopentyl | CF ₃ CH ₂ |

| Cmpd No | R ³ | R ⁴ |
|---------|------------------------------------------------------------------------|---------------------------------|
| 596 | 1-CH ₃ -cyclohexyl | CF ₃ CH ₂ |
| 597 | <i>cis</i> -2-CH ₃ -cyclohexyl | CF ₃ CH ₂ |
| 598 | <i>trans</i> -2-CH ₃ -cyclohexyl | CF ₃ CH ₂ |
| 599 | <i>cis</i> -3-CH ₃ -cyclohexyl | CF ₃ CH ₂ |
| 600 | <i>trans</i> -3-CH ₃ -cyclohexyl | CF ₃ CH ₂ |
| 601 | 2,2-(CH ₃) ₂ -cyclohexyl | CF ₃ CH ₂ |
| 602 | 3,3-(CH ₃) ₂ -cyclohexyl | CF ₃ CH ₂ |
| 603 | <i>cis</i> -4-CH ₃ -cyclohexyl | CF ₃ CH ₂ |
| 604 | <i>trans</i> -4-CH ₃ -cyclohexyl | CF ₃ CH ₂ |
| 605 | 4,4-(CH ₃) ₂ -cyclohexyl | CF ₃ CH ₂ |
| 606 | 4-(CH ₃) ₃ C-cyclohexyl | CF ₃ CH ₂ |
| 607 | <i>cis</i> -1-CH ₃ -2-fluoro-cyclopropyl | H |
| 608 | <i>trans</i> -1-CH ₃ -2-fluoro-cyclopropyl | H |
| 609 | 1-CH ₃ -2,2-difluoro-cyclopropyl | H |
| 610 | <i>cis</i> -1-CH ₃ -2-chloro-2-fluorocyclopropyl | H |
| 611 | <i>trans</i> -1-CH ₃ -2-chloro-2-fluorocyclopropyl | H |
| 612 | CH ₃ CO(CH ₃)CH | H |
| 613 | CH ₃ CH ₂ CO(CH ₃)CH | H |
| 614 | (CH ₃) ₂ CHCO(CH ₃)CH | H |
| 615 | (CH ₃) ₃ CCO(CH ₃)CH | H |
| 616 | CH ₃ CH ₂ CH ₂ CO(CH ₃)CH | H |
| 617 | CH ₃ CO(CF ₃)CH | H |
| 618 | CH ₃ CH ₂ CO(CF ₃)CH | H |
| 619 | CH ₃ CO(CH ₃) ₂ C | H |
| 620 | CH ₃ CH ₂ CO(CH ₃) ₂ C | H |
| 621 | <i>cis</i> -1-CH ₃ -2-fluoro-cyclopropyl | CH ₃ |
| 622 | <i>trans</i> -1-CH ₃ -2-fluoro-cyclopropyl | CH ₃ |
| 623 | 1-CH ₃ -2,2-difluoro-cyclopropyl | CH ₃ |
| 624 | <i>cis</i> -1-CH ₃ -2-chloro-2-fluorocyclopropyl | CH ₃ |
| 625 | <i>trans</i> -1-CH ₃ -2-chloro-2-fluorocyclopropyl | CH ₃ |
| 626 | CH ₃ CO(CH ₃)CH | CH ₃ |
| 627 | CH ₃ CH ₂ CO(CH ₃)CH | CH ₃ |
| 628 | (CH ₃) ₂ CHCO(CH ₃)CH | CH ₃ |

| Cmpd No | R ³ | R ⁴ |
|---------|------------------------------------------------------------------------|-------------------------------|
| 629 | (CH ₃) ₃ CCO(CH ₃)CH | CH ₃ |
| 630 | CH ₃ CH ₂ CH ₂ CO(CH ₃)CH | CH ₃ |
| 631 | CH ₃ CO(CF ₃)CH | CH ₃ |
| 632 | CH ₃ CH ₂ CO(CF ₃)CH | CH ₃ |
| 633 | CH ₃ CO(CH ₃) ₂ C | CH ₃ |
| 634 | CH ₃ CH ₂ CO(CH ₃) ₂ C | CH ₃ |
| 635 | <i>cis</i> -1-CH ₃ -2-fluoro-cyclopropyl | C ₂ H ₅ |
| 636 | <i>trans</i> -1-CH ₃ -2-fluoro-cyclopropyl | C ₂ H ₅ |
| 637 | 1-CH ₃ -2,2-difluoro-cyclopropyl | C ₂ H ₅ |
| 638 | <i>cis</i> -1-CH ₃ -2-chloro-2-fluorocyclopropyl | C ₂ H ₅ |
| 639 | <i>trans</i> -1-CH ₃ -2-chloro-2-fluorocyclopropyl | C ₂ H ₅ |
| 640 | CH ₃ CO(CH ₃)CH | C ₂ H ₅ |
| 641 | CH ₃ CH ₂ CO(CH ₃)CH | C ₂ H ₅ |
| 642 | (CH ₃) ₂ CHCO(CH ₃)CH | C ₂ H ₅ |
| 643 | (CH ₃) ₃ CCO(CH ₃)CH | C ₂ H ₅ |
| 644 | CH ₃ CH ₂ CH ₂ CO(CH ₃)CH | C ₂ H ₅ |
| 645 | CH ₃ CO(CF ₃)CH | C ₂ H ₅ |

| Cmpd No | R ³ | R ⁴ |
|---------|------------------------------------------------------------------------|---------------------------------|
| 646 | CH ₃ CH ₂ CO(CF ₃)CH | C ₂ H ₅ |
| 647 | CH ₃ CO(CH ₃) ₂ C | C ₂ H ₅ |
| 648 | CH ₃ CH ₂ CO(CH ₃) ₂ C | C ₂ H ₅ |
| 649 | <i>cis</i> -1-CH ₃ -2-fluoro-cyclopropyl | CF ₃ CH ₂ |
| 650 | <i>trans</i> -1-CH ₃ -2-fluoro-cyclopropyl | CF ₃ CH ₂ |
| 651 | 1-CH ₃ -2,2-difluoro-cyclopropyl | CF ₃ CH ₂ |
| 652 | <i>cis</i> -1-CH ₃ -2-chloro-2-fluorocyclopropyl | CF ₃ CH ₂ |
| 653 | <i>trans</i> -1-CH ₃ -2-chloro-2-fluorocyclopropyl | CF ₃ CH ₂ |
| 654 | CH ₃ CO(CH ₃)CH | CF ₃ CH ₂ |
| 655 | CH ₃ CH ₂ CO(CH ₃)CH | CF ₃ CH ₂ |
| 656 | (CH ₃) ₂ CHCO(CH ₃)CH | CF ₃ CH ₂ |
| 657 | (CH ₃) ₃ CCO(CH ₃)CH | CF ₃ CH ₂ |
| 658 | CH ₃ CH ₂ CH ₂ CO(CH ₃)CH | CF ₃ CH ₂ |
| 659 | CH ₃ CO(CF ₃)CH | CF ₃ CH ₂ |
| 660 | CH ₃ CH ₂ CO(CF ₃)CH | CF ₃ CH ₂ |
| 661 | CH ₃ CO(CH ₃) ₂ C | CF ₃ CH ₂ |
| 662 | CH ₃ CH ₂ CO(CH ₃) ₂ C | CF ₃ CH ₂ |

Table 2

Table 2 consists of 662 compounds of the general formula (1A), where W, X and Z are N and Y is CH, R is Cl, R¹ is 2,5,6-trifluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 2 is the same as compound 1 of Table 1 except that in compound 1 of Table 2, R¹ is 2,5,6-trifluorophenyl. Similarly, compounds 2 to 662 of Table 2 are the same as compounds 2 to 662 of Table 1 except that in the compounds of Table 2, R¹ is 2,5,6-trifluorophenyl.

Table 3

Table 3 consists of 662 compounds of the general formula (1A), where W, X and Z are N and Y is CH, R is Cl, R¹ is 2,3,4,5,6-pentafluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 3 is the same as compound 1 of Table 1 except that in compound 1 of Table 3, R¹ is 2,3,4,5,6-pentafluorophenyl. Similarly, compounds 2 to 662 of Table 3 are the same as compounds 2 to 662 of Table 1 except that in the compounds of Table 3, R¹ is 2,3,4,5,6-pentafluorophenyl.

Table 4

Table 4 consists of 662 compounds of the general formula (1A), where W, X and Z are N and Y is CH, R is Cl, R¹ is 2,6-difluoro-4-methoxyphenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 4 is the same as compound 1 of Table 1 except that in compound 1 of Table 4, R¹ is 2,6-difluoro-4-methoxyphenyl. Similarly, compounds 2 to 662 of Table 4 are the same as compounds 2 to 662 of Table 1 except that in the compounds of Table 4, R¹ is 2,6-difluoro-4-methoxyphenyl.

Table 5

Table 5 consists of 662 compounds of the general formula (1A), where W, X and Z are N and Y is CH, R is Cl, R¹ is 2-fluoro-6-chlorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 5 is the same as compound 1 of Table 1 except that in compound 1 of Table 5, R¹ is 2-fluoro-6-chlorophenyl. Similarly, compounds 2 to 662 of Table 5 are the same as compounds 2 to 662 of Table 1 except that in the compounds of Table 5, R¹ is 2-fluoro-6-chlorophenyl.

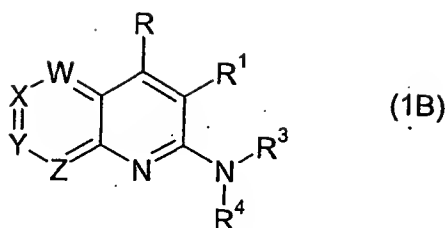
Table 6

Table 6 consists of 662 compounds of the general formula (1B), where W, X and Z are N and Y is CH, R is Cl, R¹ is 2,4,6-trifluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 6 is the same as compound 1 of Table 1 except that in compound 1 of Table 6, the compound has the general formula (1B). Similarly, compounds 2 to 662 of Table 6 are the same as compounds 2 to 662 of Table 1 except that in the compounds of Table 6, the compounds have the general formula (1B).

Table 7

Table 7 consists of 662 compounds of the general formula (1B), where W, X and Z are N and Y is CH, R is Cl, R¹ is 2,5,6-trifluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 7 is the same as compound 1 of Table 2 except that in compound 1 of Table 7, the compound has the general formula (1B). Similarly, compounds 2 to 662 of Table 7 are the same as compounds 2 to 662 of Table 2 except that in the compounds of Table 7, the compounds have the general formula (1B).

Table 8

Table 8 consists of 662 compounds of the general formula (1B), where W, X and Z are N and Y is CH, R is Cl, R¹ is 2,3,4,5,6-pentafluorophenyl, and the values of R³ and R⁴ are as listed in Table 3. Thus, compound 1 of Table 8 is the same as compound 1 of Table 3 except that in compound 1 of Table 8, the compound has the general formula (1B). Similarly, compounds 2 to 662 of Table 8 are the same as compounds 2 to 662 of Table 3 except that in the compounds of Table 8, the compounds have the general formula (1B).

Table 9

Table 9 consists of 662 compounds of the general formula (1B), where W, X and Z are N and Y is CH, R is Cl, R¹ is 2,6-difluoro-4-methoxyphenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 9 is the same as compound 1 of Table 4 except that in compound 1 of Table 9, the compound has the general formula (1B). Similarly, compounds 2 to 662 of Table 9 are the same as compounds 2 to 662 of Table 4 except that in the compounds of Table 9, the compounds have the general formula (1B).

Table 10

Table 10 consists of 662 compounds of the general formula (1B), where W, X and Z are N and Y is CH, R is Cl, R¹ is 2-fluoro-6-chlorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 10 is the same as compound 1 of Table 5 except that in compound 1 of Table 10, the compound has the general formula (1B). Similarly, compounds 2 to 662 of Table 10 are the same as compounds 2 to 662 of Table 5 except that in the compounds of Table 10, the compounds have the general formula (1B).

Table 11

Table 11 consists of 662 compounds of the general formula (1A), where W, Y and Z are N and X is CH, R is Cl, R¹ is 2,4,6-trifluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 11 is the same as compound 1 of Table 1 except that in compound 1 of Table 11, the compound has the general formula (1A) where W, Y and Z are N and X is CH. Similarly, compounds 2 to 662 of Table 11 are the same as compounds 2 to 662 of Table 1 except that in the compounds of Table 11, the compounds have the general formula (1A) where W, Y and Z are N and X is CH.

Table 12

Table 12 consists of 662 compounds of the general formula (1A), where W, Y and Z are N and X is CH, R is Cl, R¹ is 2,5,6-trifluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 12 is the same as compound 1 of Table 2 except that in compound 1 of Table 12, the compound has the general formula (1A) where W, Y and Z are N and X is CH. Similarly, compounds 2 to 662 of Table 12 are the same as compounds 2 to 662 of Table 2 except that in the compounds of Table 12, the compounds have the general formula (1A) where W, Y and Z are N and X is CH.

Table 13

Table 13 consists of 662 compounds of the general formula (1A), where W, Y and Z are N and X is CH, R is Cl, R¹ is 2,3,4,5,6-pentafluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 13 is the same as compound 1 of Table 3 except that in compound 1 of Table 13, the compound has the general formula (1A) where W, Y and Z are N and X is CH. Similarly, compounds 2 to 662 of Table 13 are the same as compounds 2 to 662 of Table 3 except that in the compounds of Table 13, the compounds have the general formula (1A) where W, Y and Z are N and X is CH.

Table 14

Table 14 consists of 662 compounds of the general formula (1A), where W, Y and Z are N and X is CH, R is Cl, R¹ is 2,6-difluoro-4-methoxyphenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 14 is the same as compound 1 of Table 4 except that in compound 1 of Table 14, the compound has the general formula (1A) where W, Y and Z are N and X is CH. Similarly, compounds 2 to 662 of Table 14 are the same as compounds 2 to 662 of Table 4 except that in the compounds of Table 14, the compounds have the general formula (1A) where W, Y and Z are N and X is CH.

Table 15

Table 15 consists of 662 compounds of the general formula (1A), where W, Y and Z are N and X is CH, R is Cl, R¹ is 2-fluoro-6-chlorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 15 is the same as compound 1 of Table 5 except that in compound 1 of Table 15, the compound has the general formula (1A) where W, Y and Z are N and X is CH. Similarly, compounds 2 to 662 of Table 15 are the same as compounds 2 to 662 of Table 5 except that in the compounds of Table 15, the compounds have the general formula (1A) where W, Y and Z are N and X is CH.

Table 16

Table 16 consists of 662 compounds of the general formula (1B), where W, Y and Z are N and X is CH, R is Cl, R¹ is 2,4,6-trifluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 16 is the same as compound 1 of Table 11 except that in compound 1 of Table 16, the compound has the general formula (1B). Similarly, compounds 2 to 662 of Table 16 are the same as compounds 2 to 662 of Table 11 except that in the compounds of Table 16, the compounds have the general formula (1B).

Table 17

Table 17 consists of 662 compounds of the general formula (1B), where W, Y and Z are N and X is CH, R is Cl, R¹ is 2,5,6-trifluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 17 is the same as compound 1 of Table 12 except that in compound 1 of Table 17, the compound has the general formula (1B). Similarly, compounds 2 to 662 of Table 17 are the same as compounds 2 to 662 of Table 12 except that in the compounds of Table 17, the compounds have the general formula (1B).

Table 18

Table 18 consists of 662 compounds of the general formula (1B), where W, Y and Z are N and X is CH, R is Cl, R¹ is 2,3,4,5,6-pentafluorophenyl, and the values of R³ and R⁴

are as listed in Table 1. Thus, compound 1 of Table 18 is the same as compound 1 of Table 13 except that in compound 1 of Table 18, the compound has the general formula (1B). Similarly, compounds 2 to 662 of Table 18 are the same as compounds 2 to 662 of Table 13 except that in the compounds of Table 18, the compounds have the general formula (1B).

5 Table 19

Table 19 consists of 662 compounds of the general formula (1B), where W, Y and Z are N and X is CH, R is Cl, R¹ is 2,6-difluoro-4-methoxyphenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 19 is the same as compound 1 of Table 14 except that in compound 1 of Table 19, the compound has the general formula (1B).

10 Similarly, compounds 2 to 662 of Table 19 are the same as compounds 2 to 662 of Table 14 except that in the compounds of Table 19, the compounds have the general formula (1B).

Table 20

Table 20 consists of 662 compounds of the general formula (1B), where W, Y and Z are N and X is CH, R is Cl, R¹ is 2-fluoro-6-chlorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 20 is the same as compound 1 of Table 15 except that in compound 1 of Table 20, the compound has the general formula (1B).

Similarly, compounds 2 to 662 of Table 20 are the same as compounds 2 to 662 of Table 15 except that in the compounds of Table 20, the compounds have the general formula (1B).

Table 21

20 Table 21 consists of 662 compounds of the general formula (1A), where W, X and Z are N and Y is CH, R is F, R¹ is 2,4,6-trifluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 21 is the same as compound 1 of Table 1 except that in compound 1 of Table 21, R is F instead of Cl. Similarly, compounds 2 to 662 of Table 21 are the same as compounds 2 to 662 of Table 1 except that in the compounds of Table 21, R is F instead of Cl.

Table 22

Table 22 consists of 662 compounds of the general formula (1A), where W, X and Z are N and Y is CH, R is F, R¹ is 2,5,6-trifluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 22 is the same as compound 1 of Table 2 except that in compound 1 of Table 22, R is F instead of Cl. Similarly, compounds 2 to 662 of Table 22 are the same as compounds 2 to 662 of Table 2 except that in the compounds of Table 22, R is F instead of Cl.

Table 23

Table 23 consists of 662 compounds of the general formula (1A), where W, X and Z are N and Y is CH, R is F, R¹ is 2,3,4,5,6-pentafluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 23 is the same as compound 1 of Table 3 except that in compound 1 of Table 23, R is F instead of Cl. Similarly, compounds 2 to 662 of Table 23 are the same as compounds 2 to 662 of Table 3 except that in the compounds of Table 23, R is F instead of Cl.

Table 24

Table 24 consists of 662 compounds of the general formula (1A), where W, X and Z are N and Y is CH, R is F, R¹ is 2,6-difluoro-4-methoxyphenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 24 is the same as compound 1 of Table 4 except that in compound 1 of Table 24, R is F instead of Cl. Similarly, compounds 2 to 662 of Table 24 are the same as compounds 2 to 662 of Table 4 except that in the compounds of Table 24, R is F instead of Cl.

Table 25

Table 25 consists of 662 compounds of the general formula (1A), where W, X and Z are N and Y is CH, R is F, R¹ is 2-fluoro-6-chlorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 25 is the same as compound 1 of Table 5 except that in compound 1 of Table 25, R is F instead of Cl. Similarly, compounds 2 to 662 of Table 25 are the same as compounds 2 to 662 of Table 5 except that in the compounds of Table 25, R is F instead of Cl.

Table 26

Table 26 consists of 662 compounds of the general formula (1A), where W, Y and Z are N and X is CH, R is F, R¹ is 2,4,6-trifluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 26 is the same as compound 1 of Table 11 except that in compound 1 of Table 26, R is F instead of Cl. Similarly, compounds 2 to 662 of Table 26 are the same as compounds 2 to 662 of Table 11 except that in the compounds of Table 26, R is F instead of Cl.

Table 27

Table 27 consists of 662 compounds of the general formula (1A), where W, Y and Z are N and X is CH, R is F, R¹ is 2,5,6-trifluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 27 is the same as compound 1 of Table 12

except that in compound 1 of Table 27, R is F instead of Cl. Similarly, compounds 2 to 662 of Table 27 are the same as compounds 2 to 662 of Table 12 except that in the compounds of Table 27, R is F instead of Cl.

Table 28

5 Table 28 consists of 662 compounds of the general formula (1A), where W, Y and Z are N and X is CH, R is F, R¹ is 2,3,4,5,6-pentafluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 28 is the same as compound 1 of Table 13 except that in compound 1 of Table 28, R is F instead of Cl. Similarly, compounds 2 to 662 of Table 28 are the same as compounds 2 to 662 of Table 13 except that in the compounds of
10 Table 28, R is F instead of Cl.

Table 29

 Table 29 consists of 662 compounds of the general formula (1A), where W, Y and Z are N and X is CH, R is F, R¹ is 2,6-difluoro-4-methoxyphenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 29 is the same as compound 1 of Table
15 14 except that in compound 1 of Table 29, R is F instead of Cl. Similarly, compounds 2 to 662 of Table 29 are the same as compounds 2 to 662 of Table 14 except that in the compounds of Table 29, R is F instead of Cl.

Table 30

 Table 30 consists of 662 compounds of the general formula (1A), where W, Y and Z
20 are N and X is CH, R is F, R¹ is 2-fluoro-6-chlorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 30 is the same as compound 1 of Table 15 except that in compound 1 of Table 30, R is F instead of Cl. Similarly, compounds 2 to 662 of Table 30 are the same as compounds 2 to 662 of Table 15 except that in the compounds of Table 30, R is F instead of Cl.

Table 31

25 Table 31 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the
30 same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are

exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 31 R¹ is 2,6-difluorophenyl instead of 2-fluoro-6-chlorophenyl.

Table 32

Table 32 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 32 R¹ is 2-fluorophenyl instead of 2-fluoro-6-chlorophenyl.

Table 33

Table 33 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 33 R¹ is 2,3,5,6-tetrafluorophenyl instead of 2-fluoro-6-chlorophenyl.

Table 34

Table 34 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 34 R¹ is 2-chloro-4,6-difluorophenyl instead of 2-fluoro-6-chlorophenyl.

Table 35

Table 35 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 35 R¹ is 2-chlorophenyl instead of 2-fluoro-6-chlorophenyl.

10 Table 36

Table 36 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 36 R¹ is 2,6-dichlorophenyl instead of 2-fluoro-6-chlorophenyl.

Table 37

20 Table 37 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 37 R¹ is 2,4-dichlorophenyl instead of 2-fluoro-6-chlorophenyl.

Table 38

30 Table 38 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the

same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 38 R¹ is 2,4,6-trichlorophenyl instead of 2-fluoro-6-chlorophenyl.

5 Table 39

Table 39 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the
10 same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 39 R¹ is 2,3,6-trichlorophenyl instead of 2-fluoro-6-chlorophenyl.

Table 40

15 Table 40 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly
20 the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 40 R¹ is pentachlorophenyl instead of 2-fluoro-6-chlorophenyl.

Table 41

Table 41 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as
25 compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are
30 exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 41 R¹ is 2-fluoro-4,6-dichlorophenyl instead of 2-fluoro-6-chlorophenyl.

Table 42

Table 42 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 42 R¹ is 4-fluoro-2,6-dichlorophenyl instead of 2-fluoro-6-chlorophenyl.

Table 43

Table 43 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 43 R¹ is 2-bromophenyl instead of 2-fluoro-6-chlorophenyl.

Table 44

Table 44 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 44 R¹ is 2-fluoro-6-bromophenyl instead of 2-fluoro-6-chlorophenyl.

Table 45

Table 45 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same

as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 45 R¹ is 2-bromo-4,6-difluorophenyl instead of 2-fluoro-6-chlorophenyl.

Table 46

Table 46 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 46 R¹ is 2-fluoro-6-methylphenyl instead of 2-fluoro-6-chlorophenyl.

Table 47

Table 47 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 47 R¹ is 2-chloro-6-methylphenyl instead of 2-fluoro-6-chlorophenyl.

Table 48

Table 48 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are

exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 48 R¹ is 2-methoxyphenyl instead of 2-fluoro-6-chlorophenyl.

Table 49

Table 49 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as
5 compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as
compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same
as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the
same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly
the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are
10 exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the
compounds of Table 49 R¹ is 2,6-dimethoxyphenyl instead of 2-fluoro-6-chlorophenyl.

Table 50

Table 50 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as
compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as
15 compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same
as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the
same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly
the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are
exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the
20 compounds of Table 50 R¹ is 2-fluoro-6-methoxyphenyl instead of 2-fluoro-6-chlorophenyl.

Table 51

Table 51 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as
compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as
compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same
25 as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the
same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly
the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are
exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the
compounds of Table 51 R¹ is 2-trifluoromethylphenyl instead of 2-fluoro-6-chlorophenyl.

30 Table 52

Table 52 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as
compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as

compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 52 R¹ is 2-fluoro-6-trifluoromethylphenyl instead of 2-fluoro-6-chlorophenyl.

Table 53

Table 53 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 53 R¹ is 2,6-di-(trifluoromethyl)phenyl instead of 2-fluoro-6-chlorophenyl.

Table 54

Table 54 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 54 R¹ is 2-chloro-6-trifluoromethylphenyl instead of 2-fluoro-6-chlorophenyl.

Table 55

Table 55 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the

same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 55 R¹ is 2,4-difluoro-6-trifluoromethylphenyl instead of 2-fluoro-6-chlorophenyl.

Table 56

Table 56 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 56 R¹ is 2,4-difluoro-6-methoxyphenyl instead of 2-fluoro-6-chlorophenyl.

Table 57

Table 57 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 57 R¹ is 2,4-difluoro-6-methylphenyl instead of 2-fluoro-6-chlorophenyl.

Table 58

Table 58 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are

exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 58 R¹ is 2,4-difluoropyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 59

Table 59 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 59 R¹ is 3,5-difluoropyrid-4-yl instead of 2-fluoro-6-chlorophenyl.

Table 60

Table 60 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 60 R¹ is tetrafluoropyrid-4-yl instead of 2-fluoro-6-chlorophenyl.

Table 61

Table 61 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 61 R¹ is 3-fluoropyrid-2-yl instead of 2-fluoro-6-chlorophenyl.

Table 62

Table 62 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as

compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 62 R¹ is 4-fluoropyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 63

Table 63 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 63 R¹ is 3-fluoropyrid-4-yl instead of 2-fluoro-6-chlorophenyl.

Table 64

Table 64 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 64 R¹ is 2-fluoropyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 65

Table 65 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are

exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 65 R¹ is 2,4,6-trifluoropyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 66

Table 66 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 66 R¹ is 3,5-difluoropyrid-2-yl instead of 2-fluoro-6-chlorophenyl.

Table 67

Table 67 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 67 R¹ is 2,6-difluoropyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 68

Table 68 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 68 R¹ is 2,4-difluoro-6-methoxypyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 69

Table 69 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 69 R¹ is 2-fluoro-4-chloropyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 70

Table 70 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 70 R¹ is 3-fluoro-5-chloropyrid-4-yl instead of 2-fluoro-6-chlorophenyl.

Table 71

Table 71 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 71 R¹ is 2-chloro-4-fluoropyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 72

Table 72 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the

same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 72 R¹ is 2,4-dichloropyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

5 Table 73

Table 73 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the
10 same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 73 R¹ is 3-chloropyrid-2-yl instead of 2-fluoro-6-chlorophenyl.

Table 74

15 Table 74 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly
20 the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 74 R¹ is 4-chloropyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 75

Table 75 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as
25 compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are
30 exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 75 R¹ is 3-chloropyrid-4-yl instead of 2-fluoro-6-chlorophenyl.

Table 76

Table 76 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 76 R¹ is 2-chloropyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

10 Table 77

Table 77 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 77 R¹ is 3-trifluoromethylpyrid-2-yl instead of 2-fluoro-6-chlorophenyl.

Table 78

20 Table 78 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 78 R¹ is 4-trifluoromethylpyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 79

30 Table 79 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the

same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 79 R¹ is 3,5-dichloropyrid-2-yl instead of 2-fluoro-6-chlorophenyl.

5 Table 80

Table 80 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the
10 same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 80 R¹ is 4,6-dichloropyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 81

15 Table 81 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly
20 the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 81 R¹ is 3-trifluoromethylpyrid-4-yl instead of 2-fluoro-6-chlorophenyl.

Table 82

Table 82 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as
25 compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are
30 exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 82 R¹ is 2-trifluoromethylpyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 83

Table 83 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 83 R¹ is 2-fluoro-4-trifluoromethylpyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 84

Table 84 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 84 R¹ is 3-fluoro-5-trifluoromethylpyrid-4-yl instead of 2-fluoro-6-chlorophenyl.

Table 85

Table 85 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 85 R¹ is 4-fluoro-2-trifluoromethylpyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 86

Table 86 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 86 R¹ is 2,6-dichloropyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 87

Table 87 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 87 R¹ is 3,5-dichloropyrid-4-yl instead of 2-fluoro-6-chlorophenyl.

Table 88

Table 88 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 88 R¹ is 3-chloro-6-trifluoromethylpyrid-2-yl instead of 2-fluoro-6-chlorophenyl.

Table 89

Table 89 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same

as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 89 R¹ is 3-fluoro-6-trifluoromethylpyrid-2-yl instead of 2-fluoro-6-chlorophenyl.

Table 90

Table 90 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 90 R¹ is pyrid-2-yl instead of 2-fluoro-6-chlorophenyl.

Table 91

Table 91 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 91 R¹ is pyrid-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 92

Table 92 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are

exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 92 R¹ is pyrid-4-yl instead of 2-fluoro-6-chlorophenyl.

Table 93

Table 93 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 93 R¹ is 3-fluorothien-2-yl instead of 2-fluoro-6-chlorophenyl.

Table 94

Table 94 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 94 R¹ is 3-chlorothien-2-yl instead of 2-fluoro-6-chlorophenyl.

Table 95

Table 95 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 95 R¹ is 2,4-difluorothien-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 96

Table 96 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as

compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 96 R¹ is 2,4-dichlorothien-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 97

Table 97 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 97 R¹ is 2,4,5-trichlorothien-3-yl instead of 2-fluoro-6-chlorophenyl.

Table 98

Table 98 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 98 R¹ is piperidino instead of 2-fluoro-6-chlorophenyl.

Table 99

Table 99 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are

exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 99 R¹ is 2-methylpiperidino instead of 2-fluoro-6-chlorophenyl.

Table 100

Table 100 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 100 R¹ is 2,6-dimethylpiperidino instead of 2-fluoro-6-chlorophenyl.

Table 101

Table 101 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 101 R¹ is morpholino instead of 2-fluoro-6-chlorophenyl.

Table 102

Table 102 consists of 3972 compounds. Compounds 1 to 662 are exactly the same as compounds 1 to 662 of Table 5 respectively, compounds 663 to 1324 are exactly the same as compounds 1 to 662 of Table 10 respectively, compounds 1325 to 1986 are exactly the same as compounds 1 to 662 of Table 15 respectively, compounds 1987 to 2648 are exactly the same as compounds 1 to 662 of Table 20 respectively, compounds 2649 to 3310 are exactly the same as compounds 1 to 662 of Table 25 respectively, and compounds 3311 to 3972 are exactly the same as compounds 1 to 662 of Table 30 respectively, except that in all of the compounds of Table 102 R¹ is 2,6-dimethylmorpholino instead of 2-fluoro-6-chlorophenyl.

Table 103

Table 103 consists of 305,844 compounds. Each of these compounds is exactly the same as the corresponding compound in Tables 1 to 102 (thus, for example, compound 1 of

Table 103 is the same as compound 1 of Table 1, compound 663 of Table 103 is the same as compound 1 of Table 2, compound 19,861 of Table 103 is the same as compound 1 of Table 31, compound 305,844 of Table 103 is the same as compound 3,972 of Table 102) except that in all of the compounds of Table 103 R is ethyl instead of Cl or F.

5 Table 104

Table 104 consists of 305,844 compounds. Each of these compounds is exactly the same as the corresponding compound in Tables 1 to 102 (thus, for example, compound 1 of Table 104 is the same as compound 1 of Table 1, compound 663 of Table 104 is the same as compound 1 of Table 2, compound 19,861 of Table 104 is the same as compound 1 of Table 31, compound 305,844 of Table 104 is the same as compound 3,972 of Table 102) except that in all of the compounds of Table 104 R is Br instead of Cl or F.

Table 105

Table 105 consists of 305,844 compounds. Each of these compounds is exactly the same as the corresponding compound in Tables 1 to 102 (thus, for example, compound 1 of Table 105 is the same as compound 1 of Table 1, compound 663 of Table 105 is the same as compound 1 of Table 2, compound 19,861 of Table 105 is the same as compound 1 of Table 31, compound 305,844 of Table 105 is the same as compound 3,972 of Table 102) except that in all of the compounds of Table 105 R is methyl instead of Cl or F.

Table 106

20 Table 106 consists of 305,844 compounds. Each of these compounds is exactly the same as the corresponding compound in Tables 1 to 102 (thus, for example, compound 1 of Table 106 is the same as compound 1 of Table 1, compound 663 of Table 106 is the same as compound 1 of Table 2, compound 19,861 of Table 106 is the same as compound 1 of Table 31, compound 305,844 of Table 106 is the same as compound 3,972 of Table 102) except that in all of the compounds of Table 106 R is H instead of Cl or F.

Table 107

30 Table 107 consists of 305,844 compounds. Each of these compounds is exactly the same as the corresponding compound in Tables 1 to 102 (thus, for example, compound 1 of Table 107 is the same as compound 1 of Table 1, compound 663 of Table 107 is the same as compound 1 of Table 2, compound 19,861 of Table 107 is the same as compound 1 of Table 31, compound 305,844 of Table 107 is the same as compound 3,972 of Table 102) except that in all of the compounds of Table 107 R is cyano instead of Cl or F.

Table 108

Table 108 consists of 305,844 compounds. Each of these compounds is exactly the same as the corresponding compound in Tables 1 to 102 (thus, for example, compound 1 of Table 108 is the same as compound 1 of Table 1, compound 663 of Table 108 is the same as compound 1 of Table 2, compound 19,861 of Table 108 is the same as compound 1 of Table 31, compound 305,844 of Table 108 is the same as compound 3,972 of Table 102) except that in all of the compounds of Table 108 R is methoxy instead of Cl or F.

Table 109

Table 109 consists of 6620 compounds. Each of these compounds is exactly the same as the corresponding compound in Tables 1 to 10 (thus, for example, compound 1 of Table 109 is the same as compound 1 of Table 1, compound 663 of Table 109 is the same as compound 1 of Table 2, etc.) except that in all of the compounds of Table 109 Y is CF instead of CH.

Table 110

Table 110 consists of 6620 compounds. Each of these compounds is exactly the same as the corresponding compound in Tables 1 to 10 (thus, for example, compound 1 of Table 110 is the same as compound 1 of Table 1, compound 663 of Table 110 is the same as compound 1 of Table 2, etc.) except that in all of the compounds of Table 110 Y is CCl instead of CH.

Table 111

Table 111 consists of 6620 compounds. Each of these compounds is exactly the same as the corresponding compound in Tables 1 to 10 (thus, for example, compound 1 of Table 111 is the same as compound 1 of Table 1, compound 663 of Table 111 is the same as compound 1 of Table 2, etc.) except that in all of the compounds of Table 111 Y is CBr instead of CH.

Table 112

Table 112 consists of 6620 compounds. Each of these compounds is exactly the same as the corresponding compound in Tables 1 to 10 (thus, for example, compound 1 of Table 112 is the same as compound 1 of Table 1, compound 663 of Table 112 is the same as compound 1 of Table 2, etc.) except that in all of the compounds of Table 112 Y is CSCH₃ instead of CH.

Table 113

Table 113 consists of 6620 compounds. Each of these compounds is exactly the same as the corresponding compound in Tables 11 to 20 (thus, for example, compound 1 of Table 113 is the same as compound 1 of Table 11, compound 663 of Table 113 is the same as compound 1 of Table 12, etc.) except that in all of the compounds of Table 113 X is CF instead of CH.

Table 114

Table 114 consists of 6620 compounds. Each of these compounds is exactly the same as the corresponding compound in Tables 11 to 20 (thus, for example, compound 1 of Table 114 is the same as compound 1 of Table 11, compound 663 of Table 114 is the same as compound 1 of Table 12, etc.) except that in all of the compounds of Table 114 X is CCl instead of CH.

Table 115

Table 115 consists of 6620 compounds. Each of these compounds is exactly the same as the corresponding compound in Tables 11 to 20 (thus, for example, compound 1 of Table 115 is the same as compound 1 of Table 11, compound 663 of Table 115 is the same as compound 1 of Table 12, etc.) except that in all of the compounds of Table 115 X is CBr instead of CH.

Table 116

Table 116 consists of 6620 compounds. Each of these compounds is exactly the same as the corresponding compound in Tables 11 to 20 (thus, for example, compound 1 of Table 116 is the same as compound 1 of Table 11, compound 663 of Table 116 is the same as compound 1 of Table 12, etc.) except that in all of the compounds of Table 116 X is CSCH₃ instead of CH.

Table 117

Table 117 consists of 662 compounds of the general formula (1B), where W, X and Z are N and Y is CH, R is F, R¹ is 2,4,6-trifluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 117 is the same as compound 1 of Table 21 except that in compound 1 of Table 117, the compound has the general formula (1B). Similarly, compounds 2 to 662 of Table 117 are the same as compounds 2 to 662 of Table 21 except that in the compounds of Table 117, the compounds have the general formula (1B).

Table 118

Table 118 consists of 662 compounds of the general formula (1B) where W, X and Z are N and Y is CH, R is F, R¹ is 2,5,6-trifluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 118 is the same as compound 1 of Table 22 except that in compound 1 of Table 118, the compound has the general formula (1B).

Similarly, compounds 2 to 662 of Table 118 are the same as compounds 2 to 662 of Table 22 except that in the compounds of Table 118, the compounds have the general formula (1B).

Table 119

Table 119 consists of 662 compounds of the general formula (1B), where W, X and Z are N and Y is CH, R is F, R¹ is 2,3,4,5,6-pentafluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 119 is the same as compound 1 of Table 23 except that in compound 1 of Table 119, the compound has the general formula (1B).

Similarly, compounds 2 to 662 of Table 119 are the same as compounds 2 to 662 of Table 23 except that in the compounds of Table 119, the compounds have the general formula (1B).

Table 120

Table 120 consists of 662 compounds of the general formula (1B), where W, X and Z are N and Y is CH, R is F, R¹ is 2,6-difluoro-4-methoxyphenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 120 is the same as compound 1 of Table 24 except that in compound 1 of Table 120, the compound has the general formula (1B).

Similarly, compounds 2 to 662 of Table 120 are the same as compounds 2 to 662 of Table 24 except that in the compounds of Table 120, the compounds have the general formula (1B).

Table 121

Table 121 consists of 662 compounds of the general formula (1B), where W, X and Z are N and Y is CH, R is F, R¹ is 2-fluoro-6-chlorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 121 is the same as compound 1 of Table 25 except that in compound 1 of Table 121, the compound has the general formula (1B).

Similarly, compounds 2 to 662 of Table 121 are the same as compounds 2 to 662 of Table 25 except that in the compounds of Table 121, the compounds have the general formula (1B).

Table 122

Table 122 consists of 662 compounds of the general formula (1B), where W, Y and Z are N and X is CH, R is F, R¹ is 2,4,6-trifluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 122 is the same as compound 1 of Table 26 except that in compound 1 of Table 122, the compound has the general formula (1B).

Similarly, compounds 2 to 662 of Table 122 are the same as compounds 2 to 662 of Table 26 except that in the compounds of Table 122, the compounds have the general formula (1B).

Table 123

Table 123 consists of 662 compounds of the general formula (1B), where W, Y and Z are N and X is CH, R is F, R¹ is 2,4,5-trifluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 123 is the same as compound 1 of Table 27 except that in compound 1 of Table 123, the compound has the general formula (1B).

Similarly, compounds 2 to 662 of Table 123 are the same as compounds 2 to 662 of Table 27 except that in the compounds of Table 123, the compounds have the general formula (1B).

Table 124

Table 124 consists of 662 compounds of the general formula (1B), where W, Y and Z are N and X is CH, R is F, R¹ is 2,3,4,5,6-pentafluorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 124 is the same as compound 1 of Table 28 except that in compound 1 of Table 124, the compound has the general formula (1B).

Similarly, compounds 2 to 662 of Table 124 are the same as compounds 2 to 662 of Table 28 except that in the compounds of Table 124, the compounds have the general formula (1B).

Table 125

Table 125 consists of 662 compounds of the general formula (1B), where W, Y and Z are N and X is CH, R is F, R¹ is 2,6-difluoro-4-methoxyphenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 125 is the same as compound 1 of Table 29 except that in compound 1 of Table 125, the compound has the general formula (1B).

Similarly, compounds 2 to 662 of Table 125 are the same as compounds 2 to 662 of Table 29 except that in the compounds of Table 125, the compounds have the general formula (1B).

Table 126

Table 126 consists of 662 compounds of the general formula (1B), where W, Y and Z are N and X is CH, R is F, R¹ is 2-fluoro-6-chlorophenyl, and the values of R³ and R⁴ are as listed in Table 1. Thus, compound 1 of Table 126 is the same as compound 1 of Table 30 except that in compound 1 of Table 126, the compound has the general formula (1B).

Similarly, compounds 2 to 662 of Table 126 are the same as compounds 2 to 662 of Table 30 except that in the compounds of Table 126, the compounds have the general formula (1B).

Table 127

Table 127 consists of 662 compounds of the general formula (1A), where W, X and Z are N and Y is CH, R is NR³R⁴, R¹ is 2,4,6-trifluorophenyl and the values of R³ and R⁴ are

listed in Table 1. Thus, compound 1 of Table 127 is the same as compound 1 of Table 1 except that in compound 1 of Table 127, R is NR^3R^4 . Similarly, compounds 2 to 662 of Table 127 are the same as compounds 2 to 662 of Table 1 except that in the compounds of Table 127, R is NR^3R^4 . It should be noted that in each compound there are the two NR^3R^4 groups, both of which are the same. In other words R^3 and R^4 in the NR^3R^4 group shown in formula (1A) and in the NR^3R^4 group that is R, have the same values. These are the values set out in Table 1.

Table 128

Table 128 consists of 662 compounds of the general formula (1A), where W, X and Z are N and Y is CH, R is NR^3R^4 , R^1 is 2-fluoro-6-chlorophenyl and the values of R^3 and R^4 are listed in Table 1. Thus, compound 1 of Table 128 is the same as compound 1 of Table 5 except that in compound 1 of Table 128, R is NR^3R^4 . Similarly, compounds 2 to 662 of Table 128 are the same as compounds 2 to 662 of Table 5 except that in the compounds of Table 128, R is NR^3R^4 . It should be noted that in each compound there are the two NR^3R^4 groups, both of which are the same. In other words R^3 and R^4 in the NR^3R^4 group shown in formula (1A) and in the NR^3R^4 group that is R, have the same values. These are the values set out in Table 1.

Table 129

Table 129 consists of 662 compounds of the general formula (1A), where W, Y and Z are N and X is CH, R is NR^3R^4 , R^1 is 2,4,6-trifluorophenyl and the values of R^3 and R^4 are listed in Table 1. Thus, compound 1 of Table 129 is the same as compound 1 of Table 1 except that in compound 1 of Table 129, R is NR^3R^4 . Similarly, compounds 2 to 662 of Table 129 are the same as compounds 2 to 662 of Table 1 except that in the compounds of Table 129, R is NR^3R^4 . It should be noted that in each compound there are the two NR^3R^4 groups, both of which are the same. In other words R^3 and R^4 in the NR^3R^4 group shown in formula (1A) and in the NR^3R^4 group that is R, have the same values. These are the values set out in Table 1.

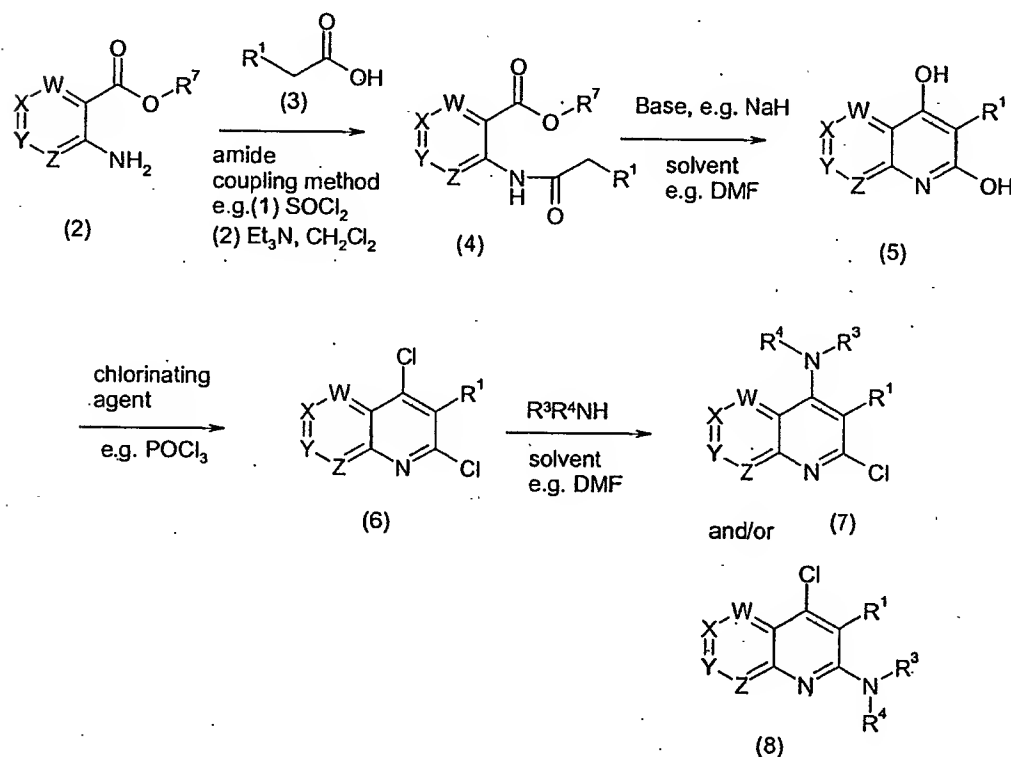
Table 130

Table 130 consists of 662 compounds of the general formula (1A), where W, Y and Z are N and X is CH, R is NR^3R^4 , R^1 is 2-fluoro-6-chlorophenyl and the values of R^3 and R^4 are listed in Table 1. Thus, compound 1 of Table 130 is the same as compound 1 of Table 5 except that in compound 1 of Table 130, R is NR^3R^4 . Similarly, compounds 2 to 662 of

Table 130 are the same as compounds 2 to 662 of Table 5 except that in the compounds of Table 130, R is NR^3R^4 . It should be noted that in each compound there are the two NR^3R^4 groups, both of which are the same. In other words R^3 and R^4 in the NR^3R^4 group shown in formula (1A) and in the NR^3R^4 group that is R, have the same values. These are the values set out in Table 1.

Compounds of formula (7) or (8), which are examples of compounds of general formula (1) where one of R and R^2 is NR^3R^4 , can be made as shown in Scheme 1, in which W, X, Y, Z, R^1 , R^3 and R^4 have the meanings given above and R^7 is C_{1-4} alkyl.

Scheme 1



10

Compounds of general formula (4) can be prepared from compounds of general formula (2), which are either commercially available or made by methods known in the literature, by reaction with acids of general formula (3), using standard coupling methods, for example by conversion to the acid chloride using a chlorinating agent such as thionyl chloride, followed by reaction of the resultant acid chloride optionally in the presence of a base such as triethylamine, in a suitable solvent such as dichloromethane or toluene.

Compounds of general formula (5) can be prepared by treating compounds of general

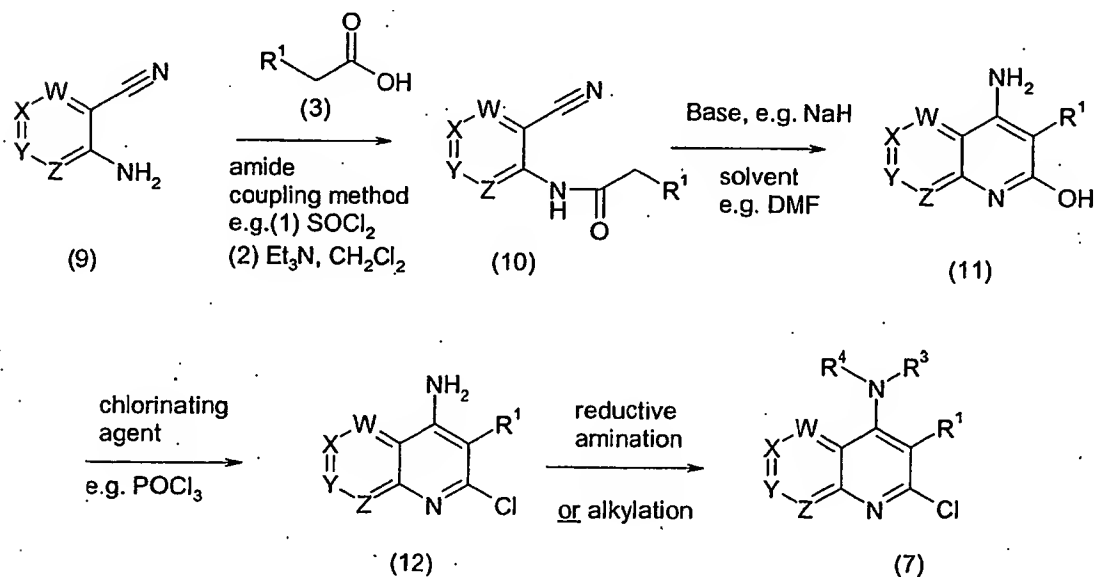
15

formula (4) with a base such as sodium hydride, optionally in the presence of a Lewis acid such as magnesium oxide, in a suitable solvent such as *N,N*-dimethylformamide (DMF) or toluene, at between room temperature and 150°C, but preferably at 60-90°C. Compounds of general formula (6) can be prepared by reaction of compounds of general formula (5) with a chlorination reagent such as phosphorus oxychloride, either neat or in a suitable solvent such as toluene, at between 50 and 150°C, but preferably between 80 and 110°C, or in a microwave reactor at between 150 and 300°C, but preferably between 200 and 250°C. Compounds of formula (7) and (8) can be prepared by reaction of compounds of general formula (6) with an amine R^3R^4NH , either neat, or in a suitable solvent such as DMF, between room temperature and 150°C, but preferably between 50 and 80°C. If compounds (7) and (8) are produced as a mixture they can be separated by suitable means such as crystallisation or chromatography under normal or reverse phase conditions.

Compounds of the general formulae (5), (6), (7) and (8) may be derivatised, via the chloro or hydroxy substituents, using routine chemical techniques to form other compounds of the general formula (1). Alternatively, other compounds of the general formula (1) may be prepared using a similar methodology to that described for preparing the compounds (5) to (8) and employing preparative techniques known from the chemical literature.

Compounds of formula (7) can also be made as shown in scheme 2.

Scheme 2



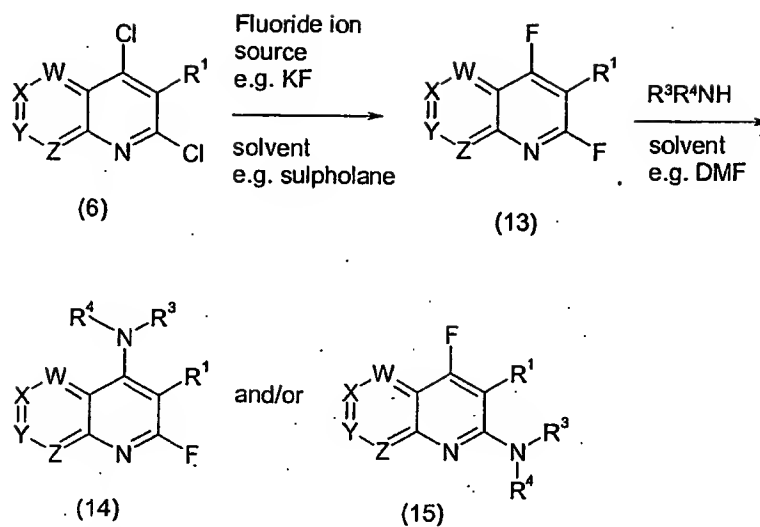
Compounds of general formula (10) can be prepared from compounds of general formula (9), which are either commercially available or made by methods known in the literature, by reaction with acids of general formula (3), using standard coupling methods, for example by conversion to the acid chloride using a chlorinating agent such as thionyl chloride, followed by reaction of the resultant acid chloride optionally in the presence of a base such as triethylamine, in a suitable solvent such as dichloromethane or toluene.

Compounds of general formula (11) can be prepared by treating compounds of general formula (10) with a base such as sodium hydride, optionally in the presence of a Lewis acid such as magnesium oxide, in a suitable solvent such as *N,N*-dimethylformamide (DMF) or toluene, at between room temperature and 150°C, but preferably at 60-90°C. Compounds of general formula (12) can be prepared by reaction of compounds of general formula (11) with a chlorination reagent such as phosphorus oxychloride, either neat or in a suitable solvent such as toluene, at between 50 and 150°C, but preferably between 80 and 110°C, or in a microwave reactor at between 150 and 300°C, but preferably between 200 and 250°C.

Compounds of formula (7) can be prepared from compounds of formula (12) by reductive amination, for example by reaction with a ketone or aldehyde in a suitable solvent such as ethanol or toluene, at between room temperature and reflux, optionally in the presence of an acid catalyst such as para-toluenesulphonic acid or a drying agent such as molecular sieves, followed by treatment with a suitable reducing agent such as sodium borohydride, at between -20°C and 40°C, but preferably at room temperature. The aldehyde or ketone is chosen so that the desired groups R³ and R⁴ are formed after reduction of the product of reaction with the amine (12). For example if compounds of formula (12) are reacted with one equivalent of propionaldehyde and then sodium borohydride, compounds of formula (7) where R³ is *n*-propyl, and R⁴ is hydrogen are formed. If required, the reaction can be repeated with a different aldehyde or ketone. For example, if acetone is used for the second reaction, then compounds of formula (7) where R³ is *n*-propyl and R⁴ is *iso*-propyl, are formed.

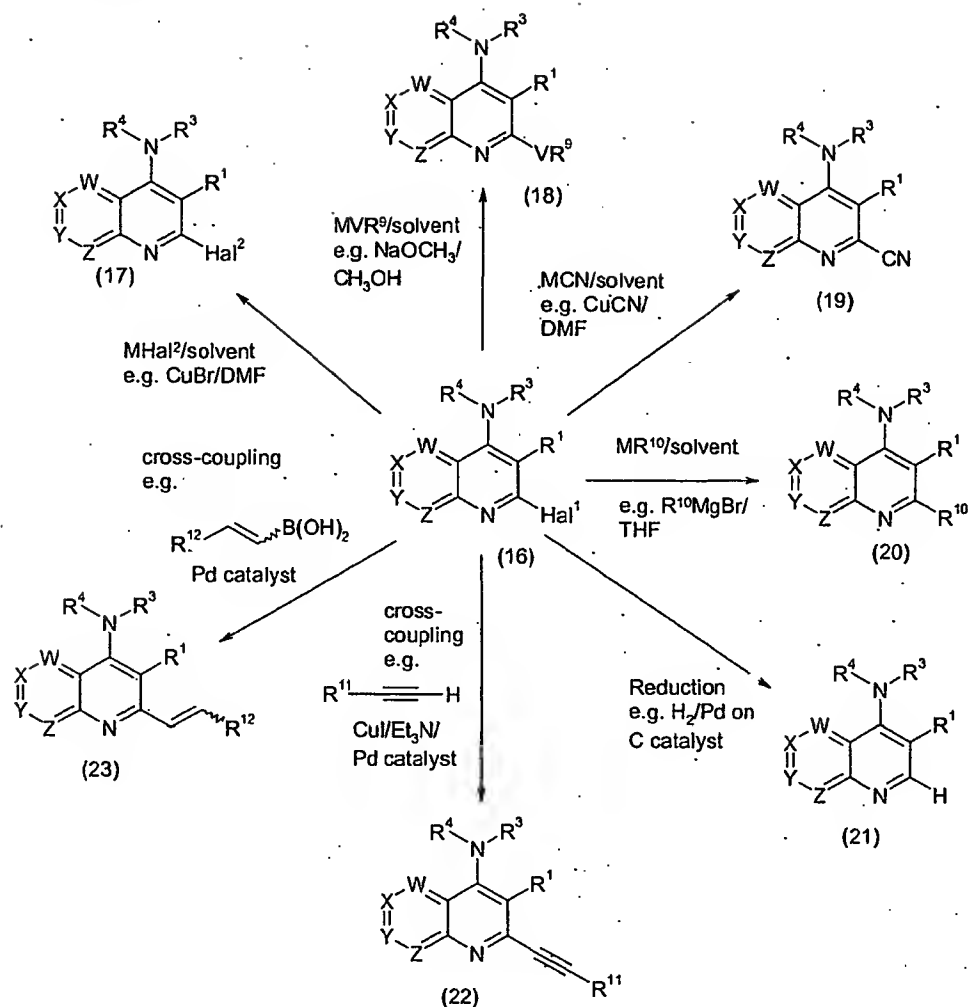
Alternatively compounds of formula (7) can be formed from compounds of formula (12) by alkylation with a group R³LG, where LG is a leaving group, by treatment with a suitable base such as sodium hydride in a solvent such as DMF, or a base such as potassium carbonate in a solvent such as acetone or DMF, at between -78°C and 100°C, but preferably between room temperature and 60°C, followed by treatment with R⁴LG in a second step under the same conditions if required.

Scheme 3



Compounds of formula (13) can be prepared as shown in Scheme 3 from compounds of formula (6) by reaction with a source of fluoride ion, such as potassium fluoride, in a suitable solvent such as sulfolane, at a temperature between 50°C and 200°C, but preferably at 80-150°C. Compounds of formula (14) and/or compounds of formula (15) can be prepared from difluoro compounds of formula (13) by reaction with an amine of formula R^3R^4NH in a suitable solvent such as DMF or CH_2Cl_2 , at a temperature of 0°C-100°C, but preferably at room temperature.

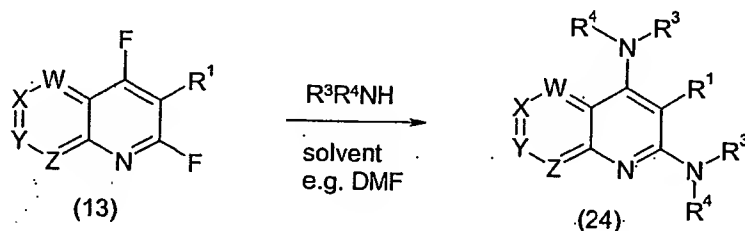
Scheme 4



Compounds of general formula (16), where Hal^1 is chlorine or fluorine, can be converted into compounds of formula (17), (18), (19), (20), (21), (22) or (23) as shown in Scheme 4. Compounds of general formula (17) where Hal^2 is bromine or iodine can be formed by reacting compounds of general formula (16) with a metal halide, for example cuprous bromide, in a suitable solvent, for example DMF, at between room temperature and 155°C, but preferably between 70°C and 155°C. Compounds of general formula (18) where V is oxygen or sulphur and R^9 is C_{1-8} alkyl, can be formed by reacting compounds of general formula (16) with a metal alkoxide or thioalkoxide MVR^9 in a suitable solvent, for example sodium methoxide in methanol, at room temperature to 65°C. Compounds of general formula (19) can be formed by reacting compounds of general formula (16) with a metal cyanide in a suitable solvent, for example cuprous cyanide in DMF, at between room temperature and

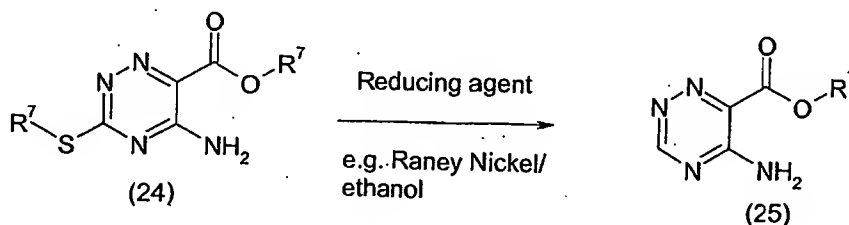
155°C but preferably between 50°C and 155°C. Compounds of general formula (20) where R^{10} is C_{1-8} alkyl, can be formed by reacting compounds of general formula (16) with an alkyl metal derivative in a suitable solvent, for example methyl magnesium bromide in THF, optionally in the presence of catalyst such as cuprous bromide or $Pd(Ph)_4$, between -40°C and 50°C. Compounds of general formula (21) can be formed by reduction of compounds of general formula (16), where Hal^1 is chlorine, for example by hydrogenolysis with hydrogen gas and a metal catalyst such as palladium on carbon in a suitable solvent such as ethanol, at room temperature. Compounds of general formula (22) where R^{11} is hydrogen or C_{1-6} alkyl, can be formed by reaction of compounds of general formula (16) with an alkyl acetylene under the Sonogashira conditions, for example with 1-propyne in triethylamine in the presence of a cuprous salt such as cuprous iodide and a palladium catalyst such as $Pd(Ph)_4$, between room temperature and 70°C. Compounds of general formula (23) where R^{12} is hydrogen or C_{1-6} alkyl, can be formed by reaction of compounds of general formula (16) with an alkenyl metal derivative in a suitable solvent, such as ethenylboronic acid in THF, in the presence of a palladium catalyst such as $Pd(Ph)_4$ and a base such as caesium carbonate, between room temperature and 65°C.

Scheme 5



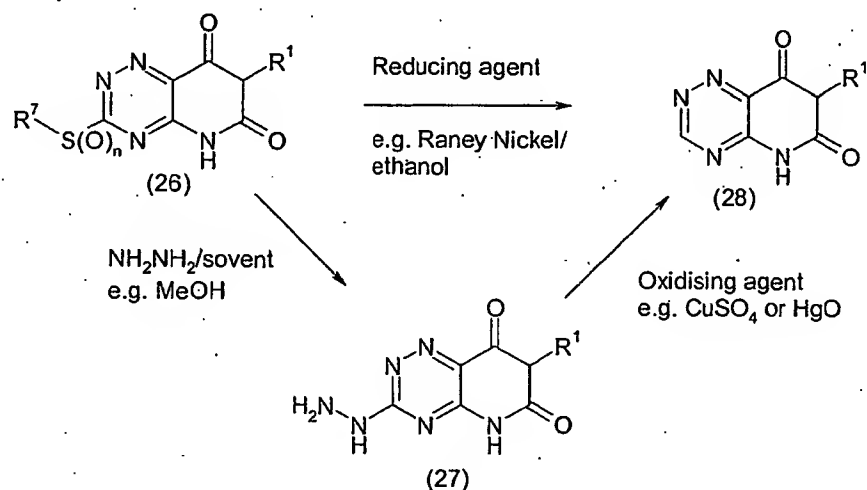
In Scheme 5 compounds of general formula (24), where the two R^3R^4N groups are identical, can be made from compounds of general formula (13) by reaction with a large excess of amine R^3R^4NH in a suitable solvent such as DMF, at a temperature between 0°C and 150°C, but preferably between room temperature and 100°C.

Scheme 6



In Scheme 6, compounds of general formula (25), which are useful for preparation as shown in Scheme 1 of compounds of formula (1) where W, X and Z are N and Y is CH, can be prepared by reduction of compounds of formula (24), which are known in the literature, using for example Raney Nickel in a suitable solvent such as ethanol, at between room temperature and reflux.

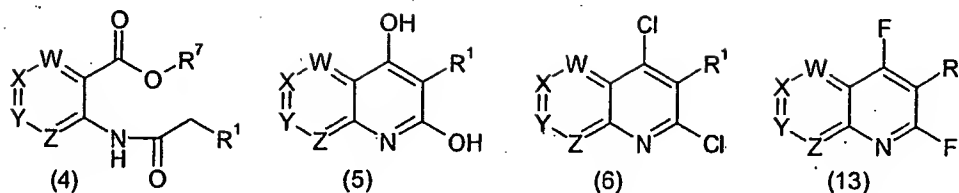
Scheme 7



In Scheme 7, compounds of general formula (28) which are useful for the preparation, as shown in Scheme 1, of compounds of formula (1) where W, X and Z are N and Y is CH, can be prepared by reduction of compounds of formula (26) where n is 0, 1 or 2, which can be prepared from compounds of formula (24) as shown in Scheme 1, using for example Raney Nickel in a suitable solvent, such as ethanol, at between room temperature and reflux. Alternatively, compounds of general formula (26) can be converted into hydrazino compounds of general formula (27), with hydrazine hydrate in a suitable solvent such as methanol, between room temperature and reflux. Compounds of formula (27) are then treated with an oxidising agent, such as copper sulphate or mercuric oxide, in a suitable solvent such as aqueous ethanol, between room temperature and reflux, to give compounds of formula (28).

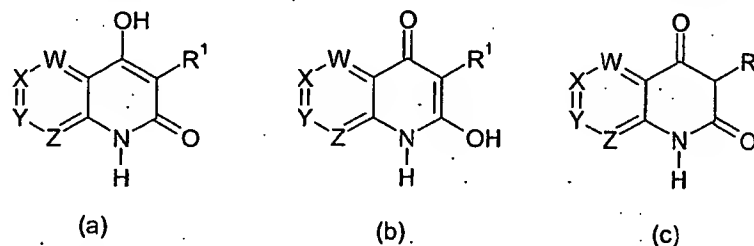
Further assistance in the preparation of the compounds of formula (1) may be derived from the following publications: Emilio, Toja, *et. al.*, *J. Heterocyclic Chem.*, 23, 1955 (1986), H. Schäfer, *et. al.*, *J. f. prakt. Chemie*, 321(4), 695 (1970) and H. Bredereck *et. al.*, *Chem. Ber.* 96, 1868-1872 (1993).

The intermediate chemicals having the general formulae (4), (5), (6) and (13):



wherein W, X, Y, Z, R¹ and R⁷ are as define above, are believed to be novel compounds and form a further part of this invention.

It should be noted that the intermediate of general formula (5) may exist in the tautomeric forms (a), (b) and (c) as well as in the form shown in formula (5):



The invention as defined by the general formula (5) embraces all such tautomers.

Of particular interest are the intermediates listed in Tables 131 to 135 below. In Table 131 the compounds have the general formula (4) where R⁷ is methyl and W, X, Y, Z and R¹ have the values shown in the table.

Table 131

| Cmpd No. | R ¹ | W | X | Y | Z |
|----------|------------------------------|---|---|----|---|
| 1 | 2,4,6-trifluorophenyl | N | N | CH | N |
| 2 | 2,5,6-trifluorophenyl | N | N | CH | N |
| 3 | 2,3,4,5,6-pentafluorophenyl | N | N | CH | N |
| 4 | 2,3,5,6-tetrafluorophenyl | N | N | CH | N |
| 5 | 2,6-difluoro-4-methoxyphenyl | N | N | CH | N |
| 6 | 2-fluoro-6-chlorophenyl | N | N | CH | N |
| 7 | 2,6-difluorophenyl | N | N | CH | N |
| 8 | 2,3,5,6-tetrafluorophenyl | N | N | CH | N |
| 9 | 2-fluorophenyl | N | N | CH | N |
| 10 | 2-chlorophenyl | N | N | CH | N |
| 11 | 2-bromophenyl | N | N | CH | N |

| | | | | | |
|----|------------------------------|---|----|-------------------|---|
| 12 | 2,4-dichlorophenyl | N | N | CH | N |
| 13 | 2,4,6-trifluorophenyl | N | CH | N | N |
| 14 | 2,5,6-trifluorophenyl | N | CH | N | N |
| 15 | 2,3,4,5,6-pentafluorophenyl | N | CH | N | N |
| 16 | 2,3,5,6-tetrafluorophenyl | N | CH | N | N |
| 17 | 2,6-difluoro-4-methoxyphenyl | N | CH | N | N |
| 18 | 2-fluoro-6-chlorophenyl | N | CH | N | N |
| 19 | 2,6-difluorophenyl | N | CH | N | N |
| 20 | 2,3,5,6-tetrafluorophenyl | N | CH | N | N |
| 21 | 2-fluorophenyl | N | CH | N | N |
| 22 | 2-chlorophenyl | N | CH | N | N |
| 23 | 2-bromophenyl | N | CH | N | N |
| 24 | 2,4-dichlorophenyl | N | CH | N | N |
| 25 | 2-fluoro-6-chlorophenyl | N | N | CSCH ₃ | N |

Table 132

Table 132 consists of 25 compounds of the general formula (5), where W, X, Y, Z and R¹ have the values given in Table 131. Thus, compound 1 of Table 132 has the same W, X, Y, Z and R¹ values as compound 1 of Table 131, etc.

Table 133

Table 133 consists of 25 compounds of the general formula (6), where W, X, Y, Z and R¹ have the values given in Table 131. Thus, compound 1 of Table 133 has the same W, X, Y, Z and R¹ values as compound 1 of Table 131, etc.

10 Table 134

Table 134 consists of 25 compounds of the general formula (13), where W, X, Y, Z and R¹ have the values given in Table 131. Thus, compound 1 of Table 134 has the same W, X, Y, Z and R¹ values as compound 1 of Table 131, etc.

Table 135

15 Table 135 consists of 25 compounds of the general formula (4), where W, X, Y, Z and R¹ have the values given in Table 131 and R⁷ is ethyl. Thus, compound 1 of Table 135 is the same as compound 1 of Table 131 except that in compound 1 of Table 135, R⁷ is ethyl.

instead of methyl. Similarly, compounds 2 to 24 of Table 135 are the same as compounds 2 to 24 of Table 131 except that in the compounds of Table 135, R⁷ is ethyl.

The compounds of formula (1) are active fungicides and may be used to control one or more of the following pathogens: *Pyricularia oryzae* (*Magnaporthe grisea*) on rice and wheat and other *Pyricularia* spp. on other hosts; *Puccinia triticina* (or *recondita*), *Puccinia striiformis* and other rusts on wheat, *Puccinia hordei*, *Puccinia striiformis* and other rusts on barley, and rusts on other hosts (for example turf, rye, coffee, pears, apples, peanuts, sugar beet, vegetables and ornamental plants); *Erysiphe cichoracearum* on cucurbits (for example melon); *Blumeria* (or *Erysiphe*) *graminis* (powdery mildew) on barley, wheat, rye and turf and other powdery mildews on various hosts, such as *Sphaerotheca macularis* on hops, *Sphaerotheca fusca* (*Sphaerotheca fuliginea*) on cucurbits (for example cucumber), *Leveillula taurica* on tomatoes, aubergine and green pepper, *Podosphaera leucotricha* on apples and *Uncinula necator* on vines; *Cochliobolus* spp., *Helminthosporium* spp., *Drechslera* spp. (*Pyrenophora* spp.), *Rhynchosporium* spp., *Mycosphaerella graminicola* (*Septoria tritici*) and *Phaeosphaeria nodorum* (*Stagonospora nodorum* or *Septoria nodorum*), *Pseudocercospora herpotrichoides* and *Gaeumannomyces graminis* on cereals (for example wheat, barley, rye), turf and other hosts; *Cercospora arachidicola* and *Cercosporidium personatum* on peanuts and other *Cercospora* spp. on other hosts, for example sugar beet, bananas, soya beans and rice; *Botrytis cinerea* (grey mould) on tomatoes, strawberries, vegetables, vines and other hosts and other *Botrytis* spp. on other hosts; *Alternaria* spp. on vegetables (for example carrots); oil-seed rape, apples, tomatoes, potatoes, cereals (for example wheat) and other hosts; *Venturia* spp. (including *Venturia inaequalis* (scab)) on apples, pears, stone fruit, tree nuts and other hosts; *Cladosporium* spp. on a range of hosts including cereals (for example wheat) and tomatoes; *Monilinia* spp. on stone fruit, tree nuts and other hosts; *Didymella* spp. on tomatoes, turf, wheat, cucurbits and other hosts; *Phoma* spp. on oil-seed rape, turf, rice, potatoes, wheat and other hosts; *Aspergillus* spp. and *Aureobasidium* spp. on wheat, lumber and other hosts; *Ascochyta* spp. on peas, wheat, barley and other hosts; *Stemphylium* spp. (*Pleospora* spp.) on apples, pears, onions and other hosts; summer diseases (for example bitter rot (*Glomerella cingulata*), black rot or frog-eye leaf spot (*Botryosphaeria obtusa*), Brooks fruit spot (*Mycosphaerella pomi*), Cedar apple rust (*Gymnosporangium juniperi-virginianae*), sooty blotch (*Gloeodes pomigena*), flyspeck (*Schizothyrium pomi*) and white rot (*Botryosphaeria dothidea*)) on apples and pears;

Plasmopara viticola on vines; other downy mildews, such as *Bremia lactucae* on lettuce, *Peronospora* spp. on soybeans, tobacco, onions and other hosts, *Pseudoperonospora humuli* on hops and *Pseudoperonospora cubensis* on cucurbits; *Pythium* spp. (including *Pythium ultimum*) on turf and other hosts; *Phytophthora infestans* on potatoes and tomatoes and other
5 *Phytophthora* spp. on vegetables, strawberries, avocado, pepper, ornamentals, tobacco, cocoa and other hosts; *Thanatephorus cucumeris* on rice and turf and other *Rhizoctonia* spp. on various hosts such as wheat and barley, peanuts, vegetables, cotton and turf; *Sclerotinia* spp. on turf, peanuts, potatoes, oil-seed rape and other hosts; *Sclerotium* spp. on turf, peanuts and other hosts; *Gibberella fujikuroi* on rice; *Colletotrichum* spp. on a range of hosts
10 including turf, coffee and vegetables; *Laetisaria fuciformis* on turf; *Mycosphaerella* spp. on bananas, peanuts, citrus, pecans, papaya and other hosts; *Diaporthe* spp. on citrus, soybean, melon, pears, lupin and other hosts; *Elsinoe* spp. on citrus, vines, olives, pecans, roses and other hosts; *Verticillium* spp. on a range of hosts including hops, potatoes and tomatoes; *Pyrenopeziza* spp. on oil-seed rape and other hosts; *Oncobasidium theobromae* on cocoa
15 causing vascular streak dieback; *Fusarium* spp., *Typhula* spp., *Microdochium nivale*, *Ustilago* spp., *Urocystis* spp., *Tilletia* spp. and *Claviceps purpurea* on a variety of hosts but particularly wheat, barley, turf and maize; *Ramularia* spp. on sugar beet, barley and other hosts; post-harvest diseases particularly of fruit (for example *Penicillium digitatum*, *Penicillium italicum* and *Trichoderma viride* on oranges, *Colletotrichum musae* and
20 *Gloeosporium musarum* on bananas and *Botrytis cinerea* on grapes); other pathogens on vines, notably *Eutypa lata*, *Guignardia bidwellii*, *Phellinus igniarius*, *Phomopsis viticola*, *Pseudopeziza tracheiphila* and *Stereum hirsutum*; other pathogens on trees (for example *Lophodermium seditiosum*) or lumber, notably *Cephaloscypha fragrans*, *Ceratocystis* spp., *Ophiostoma piceae*, *Penicillium* spp., *Trichoderma pseudokoningii*, *Trichoderma viride*,
25 *Trichoderma harzianum*, *Aspergillus niger*, *Leptographium lindbergii* and *Aureobasidium pullulans*; and fungal vectors of viral diseases (for example *Polymyxa graminis* on cereals as the vector of barley yellow mosaic virus (BYMV) and *Polymyxa betae* on sugar beet as the vector of rhizomania).

A compound of formula (1) may move acropetally, basipetally or locally in plant
30 tissue to be active against one or more fungi. Moreover, a compound of formula (1) may be volatile enough to be active in the vapour phase against one or more fungi on the plant.

The invention therefore provides a method of combating or controlling phytopathogenic fungi which comprises applying a fungicidally effective amount of a compound of formula (1), or a composition containing a compound of formula (1), to a plant, to a seed of a plant, to the locus of the plant or seed or to soil or any other plant growth medium, e.g. nutrient solution.

The term "plant" as used herein includes seedlings, bushes and trees. Furthermore, the fungicidal method of the invention includes protectant, curative, systemic, eradicant and antisporeulant treatments.

The compounds of formula (1) are preferably used for agricultural, horticultural and turfgrass purposes in the form of a composition.

In order to apply a compound of formula (1) to a plant, to a seed of a plant, to the locus of the plant or seed or to soil or any other growth medium, a compound of formula (1) is usually formulated into a composition which includes, in addition to the compound of formula (1), a suitable inert diluent or carrier and, optionally, a surface active agent (SFA). SFAs are chemicals that are able to modify the properties of an interface (for example, liquid/solid, liquid/air or liquid/liquid interfaces) by lowering the interfacial tension and thereby leading to changes in other properties (for example dispersion, emulsification and wetting). It is preferred that all compositions (both solid and liquid formulations) comprise, by weight, 0.0001 to 95%, more preferably 1 to 85%, for example 5 to 60%, of a compound of formula (1). The composition is generally used for the control of fungi such that a compound of formula (1) is applied at a rate of from 0.1g to 10kg per hectare, preferably from 1g to 6kg per hectare, more preferably from 1g to 1kg per hectare.

When used in a seed dressing, a compound of formula (1) is used at a rate of 0.0001g to 10g (for example 0.001g or 0.05g), preferably 0.005g to 10g, more preferably 0.005g to 4g, per kilogram of seed.

In another aspect the present invention provides a fungicidal composition comprising a fungicidally effective amount of a compound of formula (1) and a suitable carrier or diluent therefor.

In a still further aspect the invention provides a method of combating and controlling fungi at a locus, which comprises treating the fungi, or the locus of the fungi with a fungicidally effective amount of a composition comprising a compound of formula (1).

The compositions can be chosen from a number of formulation types, including dustable powders (DP), soluble powders (SP), water soluble granules (SG), water dispersible granules (WG), wettable powders (WP), granules (GR) (slow or fast release), soluble concentrates (SL), oil miscible liquids (OL), ultra low volume liquids (UL), emulsifiable concentrates (EC), dispersible concentrates (DC), emulsions (both oil in water (EW) and water in oil (EO)), micro-emulsions (ME), suspension concentrates (SC), aerosols, fogging/smoke formulations, capsule suspensions (CS) and seed treatment formulations. The formulation type chosen in any instance will depend upon the particular purpose envisaged and the physical, chemical and biological properties of the compound of formula (1).

Dustable powders (DP) may be prepared by mixing a compound of formula (1) with one or more solid diluents (for example natural clays, kaolin, pyrophyllite, bentonite, alumina, montmorillonite, kieselguhr, chalk, diatomaceous earths, calcium phosphates, calcium and magnesium carbonates, sulphur, lime, flours, talc and other organic and inorganic solid carriers) and mechanically grinding the mixture to a fine powder.

Soluble powders (SP) may be prepared by mixing a compound of formula (1) with one or more water-soluble inorganic salts (such as sodium bicarbonate, sodium carbonate or magnesium sulphate) or one or more water-soluble organic solids (such as a polysaccharide) and, optionally, one or more wetting agents, one or more dispersing agents or a mixture of said agents to improve water dispersibility/solubility. The mixture is then ground to a fine powder. Similar compositions may also be granulated to form water soluble granules (SG).

Wettable powders (WP) may be prepared by mixing a compound of formula (1) with one or more solid diluents or carriers, one or more wetting agents and, preferably, one or more dispersing agents and, optionally, one or more suspending agents to facilitate the dispersion in liquids. The mixture is then ground to a fine powder. Similar compositions may also be granulated to form water dispersible granules (WG).

Granules (GR) may be formed either by granulating a mixture of a compound of formula (1) and one or more powdered solid diluents or carriers, or from pre-formed blank granules by absorbing a compound of formula (1) (or a solution thereof, in a suitable agent) in a porous granular material (such as pumice, attapulgite clays, fuller's earth, kieselguhr, diatomaceous earths or ground corn cobs) or by adsorbing a compound of formula (1) (or a solution thereof, in a suitable agent) on to a hard core material (such as sands, silicates, mineral carbonates, sulphates or phosphates) and drying if necessary. Agents which are

commonly used to aid absorption or adsorption include solvents (such as aliphatic and aromatic petroleum solvents, alcohols, ethers, ketones and esters) and sticking agents (such as polyvinyl acetates, polyvinyl alcohols, dextrans, sugars and vegetable oils). One or more other additives may also be included in granules (for example an emulsifying agent, wetting agent or dispersing agent).

Dispersible Concentrates (DC) may be prepared by dissolving a compound of formula (1) in water or an organic solvent, such as a ketone, alcohol or glycol ether. These solutions may contain a surface active agent (for example to improve water dilution or prevent crystallisation in a spray tank).

Emulsifiable concentrates (EC) or oil-in-water emulsions (EW) may be prepared by dissolving a compound of formula (1) in an organic solvent (optionally containing one or more wetting agents, one or more emulsifying agents or a mixture of said agents). Suitable organic solvents for use in ECs include aromatic hydrocarbons (such as alkylbenzenes or alkylnaphthalenes, exemplified by SOLVESSO 100, SOLVESSO 150 and SOLVESSO 200; SOLVESSO is a Registered Trade Mark), ketones (such as cyclohexanone or methylcyclohexanone), alcohols (such as benzyl alcohol, furfuryl alcohol or butanol), *N*-alkylpyrrolidones (such as *N*-methylpyrrolidone or *N*-octylpyrrolidone), dimethyl amides of fatty acids (such as C₈-C₁₀ fatty acid dimethylamide) and chlorinated hydrocarbons. An EC product may spontaneously emulsify on addition to water, to produce an emulsion with sufficient stability to allow spray application through appropriate equipment. Preparation of an EW involves obtaining a compound of formula (1) either as a liquid (if it is not a liquid at room temperature, it may be melted at a reasonable temperature, typically below 70°C) or in solution (by dissolving it in an appropriate solvent) and then emulsifying the resultant liquid or solution into water containing one or more SFAs, under high shear, to produce an emulsion. Suitable solvents for use in EWs include vegetable oils, chlorinated hydrocarbons (such as chlorobenzenes), aromatic solvents (such as alkylbenzenes or alkylnaphthalenes) and other appropriate organic solvents that have a low solubility in water.

Microemulsions (ME) may be prepared by mixing water with a blend of one or more solvents with one or more SFAs, to produce spontaneously a thermodynamically stable isotropic liquid formulation. A compound of formula (1) is present initially in either the water or the solvent/SFA blend. Suitable solvents for use in MEs include those hereinbefore described for use in ECs or in EWs. An ME may be either an oil-in-water or a water-in-oil

system (which system is present may be determined by conductivity measurements) and may be suitable for mixing water-soluble and oil-soluble pesticides in the same formulation. An ME is suitable for dilution into water, either remaining as a microemulsion or forming a conventional oil-in-water emulsion.

5 Suspension concentrates (SC) may comprise aqueous or non-aqueous suspensions of finely divided insoluble solid particles of a compound of formula (1). SCs may be prepared by ball or bead milling the solid compound of formula (1) in a suitable medium, optionally with one or more dispersing agents, to produce a fine particle suspension of the compound. One or more wetting agents may be included in the composition and a suspending agent may
10 be included to reduce the rate at which the particles settle. Alternatively, a compound of formula (1) may be dry milled and added to water, containing agents hereinbefore described, to produce the desired end product.

 Aerosol formulations comprise a compound of formula (1) and a suitable propellant (for example *n*-butane). A compound of formula (1) may also be dissolved or dispersed in a
15 suitable medium (for example water or a water miscible liquid, such as *n*-propanol) to provide compositions for use in non-pressurised, hand-actuated spray pumps.

 A compound of formula (1) may be mixed in the dry state with a pyrotechnic mixture to form a composition suitable for generating, in an enclosed space, a smoke containing the compound.

20 Capsule suspensions (CS) may be prepared in a manner similar to the preparation of EW formulations but with an additional polymerisation stage such that an aqueous dispersion of oil droplets is obtained, in which each oil droplet is encapsulated by a polymeric shell and contains a compound of formula (1) and, optionally, a carrier or diluent therefor. The polymeric shell may be produced by either an interfacial polycondensation reaction or by a
25 coacervation procedure. The compositions may provide for controlled release of the compound of formula (1) and they may be used for seed treatment. A compound of formula (1) may also be formulated in a biodegradable polymeric matrix to provide a slow, controlled release of the compound.

 A composition may include one or more additives to improve the biological
30 performance of the composition (for example by improving wetting, retention or distribution on surfaces; resistance to rain on treated surfaces; or uptake or mobility of a compound of formula (1)). Such additives include surface active agents, spray additives based on oils, for

example certain mineral oils or natural plant oils (such as soy bean and rape seed oil), and blends of these with other bio-enhancing adjuvants (ingredients which may aid or modify the action of a compound of formula (1)).

A compound of formula (1) may also be formulated for use as a seed treatment, for example as a powder composition, including a powder for dry seed treatment (DS), a water soluble powder (SS) or a water dispersible powder for slurry treatment (WS), or as a liquid composition, including a flowable concentrate (FS), a solution (LS) or a capsule suspension (CS). The preparations of DS, SS, WS, FS and LS compositions are very similar to those of, respectively, DP, SP, WP, SC and DC compositions described above. Compositions for treating seed may include an agent for assisting the adhesion of the composition to the seed (for example a mineral oil or a film-forming barrier).

Wetting agents, dispersing agents and emulsifying agents may be SFAs of the cationic, anionic, amphoteric or non-ionic type.

Suitable SFAs of the cationic type include quaternary ammonium compounds (for example cetyltrimethyl ammonium bromide), imidazolines and amine salts.

Suitable anionic SFAs include alkali metals salts of fatty acids, salts of aliphatic monoesters of sulphuric acid (for example sodium lauryl sulphate), salts of sulphonated aromatic compounds (for example sodium dodecylbenzenesulphonate, calcium dodecylbenzenesulphonate, butyl-naphthalene sulphonate and mixtures of sodium di-*isopropyl*- and tri-*isopropyl*-naphthalene sulphonates), ether sulphates, alcohol ether sulphates (for example sodium laureth-3-sulphate), ether carboxylates (for example sodium laureth-3-carboxylate), phosphate esters (products from the reaction between one or more fatty alcohols and phosphoric acid (predominately mono-esters) or phosphorus pentoxide (predominately di-esters), for example the reaction between lauryl alcohol and tetraphosphoric acid; additionally these products may be ethoxylated), sulphosuccinamates, paraffin or olefine sulphonates, taurates and lignosulphonates.

Suitable SFAs of the amphoteric type include betaines, propionates and glycines.

Suitable SFAs of the non-ionic type include condensation products of alkylene oxides, such as ethylene oxide, propylene oxide, butylene oxide or mixtures thereof, with fatty alcohols (such as oleyl alcohol or cetyl alcohol) or with alkylphenols (such as octylphenol, nonylphenol or octylcresol); partial esters derived from long chain fatty acids or hexitol anhydrides; condensation products of said partial esters with ethylene oxide; block polymers

(comprising ethylene oxide and propylene oxide); alkanolamides; simple esters (for example fatty acid polyethylene glycol esters); amine oxides (for example lauryl dimethyl amine oxide); and lecithins.

Suitable suspending agents include hydrophilic colloids (such as polysaccharides, polyvinylpyrrolidone or sodium carboxymethylcellulose) and swelling clays (such as bentonite or attapulgite).

A compound of formula (1) may be applied by any of the known means of applying fungicidal compounds. For example, it may be applied, formulated or unformulated, to any part of the plant, including the foliage, stems, branches or roots, to the seed before it is planted or to other media in which plants are growing or are to be planted (such as soil surrounding the roots, the soil generally, paddy water or hydroponic culture systems), directly or it may be sprayed on, dusted on, applied by dipping, applied as a cream or paste formulation, applied as a vapour or applied through distribution or incorporation of a composition (such as a granular composition or a composition packed in a water-soluble bag) in soil or an aqueous environment.

A compound of formula (1) may also be injected into plants or sprayed onto vegetation using electrodynamic spraying techniques or other low volume methods, or applied by land or aerial irrigation systems.

Compositions for use as aqueous preparations (aqueous solutions or dispersions) are generally supplied in the form of a concentrate containing a high proportion of the active ingredient, the concentrate being added to water before use. These concentrates, which may include DCs, SCs, ECs, EWs, MEs SGs, SPs, WPs, WGs and CSs, are often required to withstand storage for prolonged periods and, after such storage, to be capable of addition to water to form aqueous preparations which remain homogeneous for a sufficient time to enable them to be applied by conventional spray equipment. Such aqueous preparations may contain varying amounts of a compound of formula (1) (for example 0.0001 to 10%, by weight) depending upon the purpose for which they are to be used.

A compound of formula (1) may be used in mixtures with fertilisers (for example nitrogen-, potassium- or phosphorus-containing fertilisers). Suitable formulation types include granules of fertiliser. The mixtures suitably contain up to 25% by weight of the compound of formula (1).

The invention therefore also provides a fertiliser composition comprising a fertiliser and a compound of formula (1).

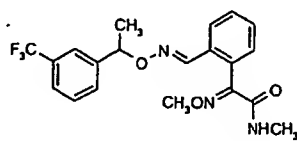
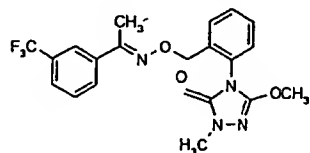
The compositions of this invention may contain other compounds having biological activity, for example micronutrients or compounds having similar or complementary
5 fungicidal activity or which possess plant growth regulating, herbicidal, insecticidal, nematicidal or acaricidal activity.

By including another fungicide, the resulting composition may have a broader spectrum of activity or a greater level of intrinsic activity than the compound of formula (1) alone. Further the other fungicide may have a synergistic effect on the fungicidal activity of
10 the compound of formula (1).

The compound of formula (1) may be the sole active ingredient of the composition or it may be admixed with one or more additional active ingredients such as a pesticide, fungicide, synergist, herbicide or plant growth regulator where appropriate. An additional active ingredient may: provide a composition having a broader spectrum of activity or
15 increased persistence at a locus; synergise the activity or complement the activity (for example by increasing the speed of effect or overcoming repellency) of the compound of formula (1); or help to overcome or prevent the development of resistance to individual components. The particular additional active ingredient will depend upon the intended utility of the composition.

20 Examples of fungicidal compounds which may be included in the composition of the invention are AC 382042 (*N*-(1-cyano-1,2-dimethylpropyl)-2-(2,4-dichlorophenoxy) propionamide), acibenzolar-S-methyl, alanycarb, aldimorph, anilazine, azaconazole, azafenidin, azoxystrobin, benalaxyl, benomyl, bentiavalicarb, biloxazol, bitertanol, blastidicin S, boscalid (new name for nicobifen), bromuconazole, bupirimate, captafol, captan,
25 carbendazim, carbendazim chlorhydrate, carboxin, carpropamid, carvone, CGA 41396, CGA 41397, chinomethionate, chlorbenzthiazole, chlorothalonil, chlorozolate, clozylacon, copper containing compounds such as copper oxychloride, copper oxyquinolate, copper sulphate, copper tallate, and Bordeaux mixture, cyamidazosulfamid, cyazofamid (IKF-916), cyflufenamid, cymoxanil, cyproconazole, cyprodinil, debacarb, di-2-pyridyl disulphide
30 1,1'-dioxide, dichlofluanid, diclocymet, diclomezine, dicloran, diethofencarb, difenoconazole, difenzoquat, diflumetorim, *O,O*-di-*iso*-propyl-*S*-benzyl thiophosphate, dimefluazole, dimetconazole, dimethirimol, dimethomorph, dimoxystrobin, diniconazole,

- dinocap, dithianon, dodecyl dimethyl ammonium chloride, dodemorph, dodine, doguadine, edifenphos, epoxiconazole, ethaboxam, ethirimol, ethyl (Z)-N-benzyl-N[(methyl(methylthioethylideneaminooxycarbonyl)amino]thio)- β -alaninate, etridiazole, famoxadone, fenamidone, fenarimol, fenbuconazole, fenfuram, fenhexamid, fenoxanil (AC 382042),
- 5 fenpiclonil, fenpropidin, fenpropimorph, fentin acetate, fentin hydroxide, ferbam, ferimzone, fluazinam, fludioxonil, flumetover, flumorph, fluoroimide, fluoxastrobin, fluquinconazole, flusilazole, flusulfamide, flutolanil, flutriafol, folpet, fosetyl-aluminium, fuberidazole, furalaxyl, furametpyr, guazatine, hexaconazole, hydroxyisoxazole, hymexazole, imazalil, imibenconazole, iminoctadine, iminoctadine triacetate, ipconazole, iprobenfos, iprodione,
- 10 iprovalicarb, isopropanyl butyl carbamate, isoprothiolane, kasugamycin, kresoxim-methyl, LY186054, LY211795, LY 248908, mancozeb, maneb, mefenoxam, mepanipyrim, mepronil, metalaxyl, metalaxyl M, metconazole, metiram, metiram-zinc, metominostrobin, metrafenone, MON65500 (N-allyl-4,5-dimethyl-2-trimethylsilylthiophene-3-carboxamide), myclobutanil, NTN0301, neoasozin, nickel dimethyldithiocarbamate, nitrothale-isopropyl,
- 15 nuarimol, ofurace, organomercury compounds, orysastrobin, oxadixyl, oxasulfuron, oxolinic acid, oxpoconazole, oxycarboxin, pefurazoate, penconazole, pencycuron, phenazin oxide, phosphorus acids, phthalide, picoxystrobin, polyoxin D, polyram, probenazole, prochloraz, procymidone, propamocarb, propamocarb hydrochloride, propiconazole, propineb, propionic acid, proquinazid, prothioconazole, pyraclostrobin, pyrazophos, pyrifenoxy, pyrimethanil,
- 20 pyroquilon, pyroxyfur, pyrrolnitrin, quaternary ammonium compounds, quinomethionate, quinoxifen, quintozone, silthiofam (MON 65500), S-imazalil, simeconazole, sipconazole, sodium pentachlorophenate, spiroxamine, streptomycin, sulphur, tebuconazole, tecloftalam, tecnazene, tetraconazole, thiabendazole, thifluzamide, 2-(thiocyanomethylthio)benzothiazole, thiophanate-methyl, thiram, tiadinil, timibenconazole, tolclofos-methyl, tolylfluanid,
- 25 triadimefon, triadimenol, triazbutil, triazoxide, tricyclazole, tridemorph, trifloxystrobin, triflumizole, triforine, triticonazole, validamycin A, vapam, vinclozolin, XRD-563, zineb, ziram, zoxamide and compounds of the formulae:



The compounds of formula (1) may be mixed with soil, peat or other rooting media for the protection of plants against seed-borne, soil-borne or foliar fungal diseases.

Some mixtures may comprise active ingredients, which have significantly different physical, chemical or biological properties such that they do not easily lend themselves to the same conventional formulation type. In these circumstances other formulation types may be prepared. For example, where one active ingredient is a water insoluble solid and the other a water insoluble liquid, it may nevertheless be possible to disperse each active ingredient in the same continuous aqueous phase by dispersing the solid active ingredient as a suspension (using a preparation analogous to that of an SC) but dispersing the liquid active ingredient as an emulsion (using a preparation analogous to that of an EW). The resultant composition is a suspoemulsion (SE) formulation.

The invention is illustrated by the following Examples in which the following abbreviations are used:

ml = millilitres

g = grammes

ppm = parts per million

s = singlet

d = doublet

t = triplet

q = quartet

m = multiplet

b = broad

f = fine

THF = tetrahydrofuran

DCM = dichloromethane

DMF = *N, N*-dimethylformamide

DMSO = dimethylsulphoxide

DMAP = 4-dimethylaminopyridine

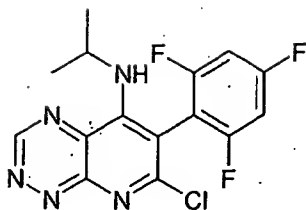
NMR = nuclear magnetic resonance

HPLC = high performance liquid

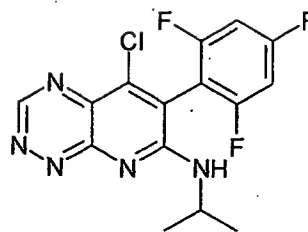
chromatography

EXAMPLE 1

- 15 This Example illustrates the preparation of [7-chloro-6-(2,4,6-trifluorophenyl)-pyrido[3,2-*e*]-[1,2,4]triazin-5-yl]-isopropylamine, Compound No.3 from Table 11, and [5-chloro-6-(2,4,6-trifluorophenyl)-pyrido[3,2-*e*][1,2,4]triazin-7-yl]-isopropylamine, Compound No.3 from Table 16.



Compound No. 3, Table 11



Compound No. 3, Table 16

Step 1

4,6-Dichloropyrimidin-5-ylamine (20.0g) was stirred while hydrazine hydrate (200ml) was added to the solid rapidly over 2 minutes. The texture of the solid changed and the mobile phase looked greenish, and the temperature rose to 35°C. The reaction was then cooled, the thick solid was collected and washed with isopropyl alcohol, and was dried to (4,6-dichloro-pyrimidin-5-yl)-hydrazine as a white fluffy solid (16.4g).

^1H NMR (d^6 -DMSO) δ ppm: 4.4-4.6 (bs, 3H), 7.75 (s, 1H), 8.15 (bs, 2H)

Step 2

The product from Step 1 (10.6g) was suspended in trimethylorthoformate (150ml), followed by the addition of concentrated hydrochloric acid (2ml). The colour changed to orange and the reaction was brought to reflux for 6 hours with stirring. The bright yellow reaction mixture was cooled and the solid collected, washed with diethyl ether and dried overnight to give 5-chloro-1,2-dihydropyrimido[5,4-e][1,2,4]triazine as a yellow solid (5.5g).

^1H NMR (d^6 -DMSO) δ ppm: 6.2 (fd, 1H), 7.4 (s, 1H), 8.0 (bs, 1H), 9.1 (bs, 1H)

Step 3

The product from Step 2 (5.5g) was suspended in ethanol (350ml) and a solution of bromine (3.1ml) in ethanol (50ml) was added slowly over a period of 10 minutes. The suspension dissolved to give a dark straw colour. After 15 minutes of stirring, water (5ml) was added and stirring was continued for 3 hours. The reaction mixture was then allowed to stand for 18 hours. The liquid was decanted from the solid, which was washed with diethyl ether, collected and discarded. The combined washings were evaporated to give a brown sludge, which was dissolved in water (5ml) and neutralised with aqueous sodium bicarbonate and then extracted with DCM (2 x 200ml) to give a yellow solution. This solution was dried over magnesium sulphate and evaporated to give 6-amino-[1,2,4]triazine-5-carboxylic acid ethyl ester as a straw solid (3.6g).

^1H NMR (CDCl_3) δ ppm: 1.5 (t, 3H), 4.5 (q, 2H), 6.4-6.7 (bs, 2H), 9.3 (s, 1H).

Step 4

The product from Step 3 (1.4g) was dissolved in dry toluene (30 ml), and DMAP (1.0g) added to give a dark brown solution. To this was added (2,4,6-trifluoro-phenyl)-acetyl chloride (1.7g) in toluene (20ml), giving a thick precipitate. The stirred suspension was heated for 5 hours at reflux, and the reaction was then cooled and the solid collected and washed with diethyl ether to yield an off white solid. The pale red filtrate was evaporated to give a viscous liquid, which was purified by column chromatography on silica gel (40-60) eluting with diethyl ether to give 6-[2-(2,4,6-trifluorophenyl)-acetylamino]-[1,2,4]triazine-5-carboxylic acid ethyl ester as a viscous oil which solidified (1.0g).

10 Step 5

The product of Step 4 (1.0g) was dissolved in DMF (20ml) and potassium carbonate (0.5g) was added and the reaction brought to 130°C for 2 hours. The reaction was cooled and the solvent evaporated to give a solid, which was triturated with diethyl ether. The solid was dissolved in water (10ml) and acidified with 2M aqueous hydrochloric acid, which precipitated a dark brown solid (0.30g). The aqueous fraction was acidified again with 2M aqueous hydrochloric acid and extracted with ethyl acetate. The organic layers were evaporated to give a brown solid, which was triturated with diethyl ether and collected to give pure 6-(2,4,6-trifluorophenyl)-8H-pyrido[3,2-e][1,2,4]triazine-5,7-dione as a pale brown solid (0.160g).

20 ¹H NMR (CD₃OD) δ ppm: 6.95(t, 2H), 9.65(s, 1H).

Step 6

The product from Step 5 (0.05g) was dissolved in phosphorus oxychloride (3.0ml) and brought to 100°C for 2 hours. The reaction was cooled and the solvent evaporated to give an oil, which was diluted with DCM (100ml) and washed with aqueous sodium bicarbonate. The DCM fraction was filtered through phase separation paper and evaporated to give 5,7-dichloro-6-(2,4,6-trifluorophenyl)-pyrido[3,2-e][1,2,4]triazine as a reddish oil, which was used without further purification.

Step 7

30 The product from Step 6 (0.02g) was dissolved in DCM (3.0ml) and then isopropylamine (1.0ml) was added, with stirring. The reaction was stirred in a sealed tube at room temperature (32°C) for 30 minutes, the volatiles were evaporated and the residue was purified by flash column chromatography on silica gel (40-60) eluting with diethyl ether to

yield an inseparable mixture of 4:6 [7-chloro-6-(2,4,6-trifluorophenyl)-pyrido[3,2-e][1,2,4]-triazin-5-yl]-isopropylamine, and [5-chloro-6-(2,4,6-trifluorophenyl)-pyrido[3,2-e][1,2,4]-triazin-7-yl]-isopropylamine, as a yellow oil (0.006g).

[7-Chloro-6-(2,4,6-trifluorophenyl)-pyrido[3,2-e][1,2,4]triazin-5-yl]-isopropylamine:

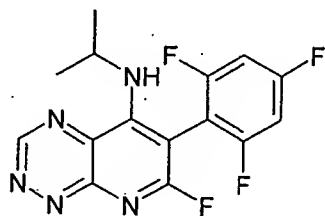
¹H NMR (CDCl₃) δ ppm: 1.15 (d,6H), 3.5 (b,1H), 6.7 (bflat,1H), 6.9 (t,2H), 9.85 (s,1H).

[5-Chloro-6-(2,4,6-trifluorophenyl)-pyrido[3,2-e][1,2,4]triazin-7-yl]-isopropylamine:

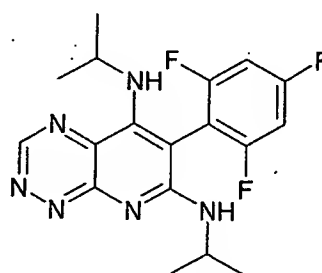
¹H NMR (CDCl₃) δ ppm: 1.3 (d,6H), 4.75 (m,1H), 4.85 (bd,1H), 7.0 (t,2H), 9.80 (s,1H).

EXAMPLE 2

This Example illustrates the preparation of [7-fluoro-6-(2,4,6-trifluorophenyl)-pyrido[3,2-e]-[1,2,4]triazin-5-yl]-isopropylamine (Compound No. 3, Table 26) and *N'*-5-,*N''*-7-diisopropyl-6-(2,4,6-trifluorophenyl)-pyrido[3,2-e][1,2,4]triazine-5,7-diamine (Compound No. 3, Table 129).



Compound No. 3, Table 26



Compound No. 3, Table 129

Step 1

5,7-Dichloro-6-(2,4,6-trifluorophenyl)-pyrido[3,2-e]triazine (1.0 g), potassium fluoride (0.53 g, spray dried) were suspended in sulpholane (5 ml) and heated to 130°C for 16 hours. After cooling, the mixture was poured into water and extracted with ethyl acetate. The organic phase was washed with water, brine and then dried over sodium sulphate. The ethyl acetate was evaporated and the residue purified by flash chromatography eluting with cyclohexane:ethyl acetate, 3:1, to give 5,7-difluoro-6-(2,4,6-trifluorophenyl)-pyrido[3,2-e]-[1,2,4]triazine as a yellow powder (0.51 g), m.p. 160-161°C.

Step 2

The product from Step 1 (0.20 g) was added to a mixture of isopropylamine (0.060g), potassium carbonate (0.140 g) and a catalytic amount of DMAP in DMF (3 ml). The mixture was stirred at room temperature for 19 hours, and then poured into ethyl acetate, washed with water and brine and the organic phase dried over sodium sulphate. The ethyl acetate was

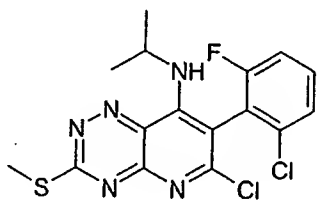
evaporated and the residue was purified by flash chromatography eluting with cyclohexane:ethyl acetate, 3:1, to give:

[7-fluoro-6-(2,4,6-trifluorophenyl)-pyrido[3,2-e][1,2,4]triazin-5-yl]-isopropylamine (0.040 g), m.p. 91-92°C, and

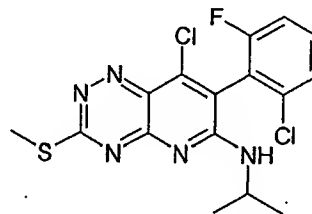
- 5 *N'*-5-,*N''*-7-diisopropyl-6-(2,4,6-trifluorophenyl)-pyrido[3,2-e][1,2,4]triazine-5,7-diamine (0.130 g), m.p. 166-167°C.

EXAMPLE 3

This Example illustrates the preparation of [6-chloro-3-methylsulfanyl-7-(2-chloro-6-fluorophenyl)-pyrido[2,3-e][1,2,4]triazin-8-yl]-isopropylamine (Compound No. 2651, Table 112) and [8-chloro-3-methylsulfanyl-7-(2-chloro-6-fluorophenyl)-pyrido[2,3-e][1,2,4]triazin-6-yl]-isopropylamine (Compound No. 5961, Table 112)



Compound No. 2651, Table 112



Compound No. 5961, Table 112

Step 1

- 5-Chloro-3-methylsulfanyl-[1,2,4]triazine-6-carboxylic acid ethyl ester (2.15 g, prepared according to the procedure in *Eur.J.Med.Chem.-Chim.Ther.* (1980), 15, 269) was dissolved in dioxane (300 ml) and cooled to 5°C. A solution of ammonia in methanol (3.0 ml of a 7M solution) was added dropwise the colour changed and a small exotherm was noticed. After about 5 minutes the reaction mixture was evaporated to give a sludge, which was extracted with ethyl acetate and was washed with water. The ethyl acetate was dried over magnesium sulphate and evaporated, and the residue purified by flash column chromatography on silica gel eluting with diethyl ether, to give 5-amino-3-methylsulfanyl-[1,2,4]-triazine-6-carboxylic acid ethyl ester as a buff solid (1.3 g).

¹H NMR (CDCl₃) δ ppm: 1.4 (t,3H), 2.6 (s,3H), 4.5 (q,2H), 5.6 (bs,1H), 7.9 (bs,1H).

Step 2

- The product from Step 1 (0.428 g) was dissolved in dry toluene (30 ml), and DMAP (0.250 g) was added to give dark a brown solution. 2-Chloro-6-fluorophenylacetyl chloride (0.414 g) in dry toluene (20 ml) was then added to give a thick precipitate, and the stirred

suspension was heated for 3 hours at reflux. The reaction was cooled and the solid collected and washed with diethyl ether to yield an off white solid. The pale red filtrate was evaporated to give a liquid that was purified by flash column chromatography eluting with diethyl ether, to give 5-[2-(2-chloro-6-fluorophenyl)-acetylamino]-3-methylsulfanyl-[1,2,4]triazine-6-
5 carboxylic acid ethyl ester as a viscous oil (0.160 g).

^1H NMR (CDCl_3) δ ppm: 1.45 (t,3H), 2.65 (s,3H), 4.5 (q,2H), 7.0 (m,1H), 7.2 (m,2H).

Step 3

The product from Step 2 (0.20 g) was dissolved in DMF (10 ml) and potassium carbonate (0.50 g) was added and the reaction brought to 130°C for 2 hours. The reaction was
10 cooled and the solvent evaporated to give a solid, which was triturated with diethyl ether. The solid was dissolved in water (10ml) and then acidified with 2M aqueous hydrochloric acid. The aqueous fraction was evaporated to give a solid, which was then triturated with methanol to give 3-methylsulfanyl-7-(2-chloro-6-fluorophenyl)-5H-pyrido[2,3-e][1,2,4]-triazine-6,8-dione as an orange solid (0.210 g).

15 ^1H NMR ($\text{D}_6\text{-DMSO}$) δ ppm: 3.5 (s,3H), 7.0 (m,1H), 7.2 (m,2H).

Step 4

The product from Step 3 (0.10 g) was suspended in 1,2-dichloroethane (5 ml) and phosphorus oxychloride (1.0 ml) was added dropwise to the suspension. The suspension dissolved, and two drops of DMF were added. The reaction was refluxed for 3 hours and
20 became dark but clear. The reaction was cooled and evaporated to give a dark brown oil. Saturated aqueous sodium bicarbonate (2 ml) was added and the mixture was extracted with DCM. The DCM fraction was dried and evaporated to give an oil, which was purified by flash column chromatography on silica gel, eluting with diethyl ether, to give 6,8-dichloro-3-methylsulfanyl-7-(2-chloro-6-fluorophenyl)-pyrido[2,3-e][1,2,4]triazine as a yellow oil
25 (0.025g).

^1H NMR (CDCl_3) δ ppm: 2.8 (s,3H), 7.35 (m,1H), 7.45 (d,1H), 7.55 (m,1H).

Step 5

To the product from Step 4 (0.025 g) in DCM (5.0 ml), isopropylamine (0.10 g) was added dropwise. The reaction was stirred at room temperature for 1 hour and the volatiles
30 were evaporated to give a yellow oil, which was purified by flash column chromatography eluting with diethyl ether, to give a 1:1 mixture of the two isomers as a yellow oil (0.015 g):

[6-chloro-3-methylsulfanyl-7-(2,4,6-trifluoro-phenyl)-pyrido[2,3-e][1,2,4]triazin-8-yl]-isopropylamine.

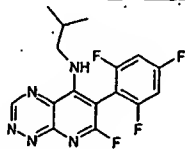
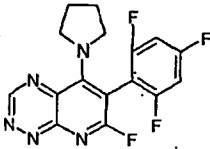
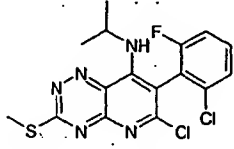
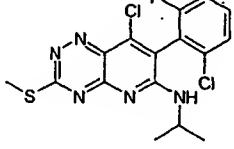
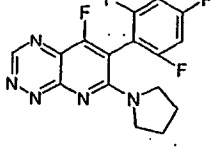
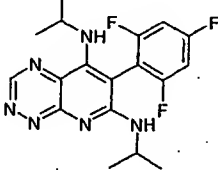
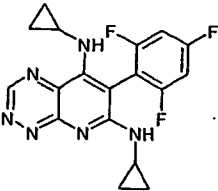
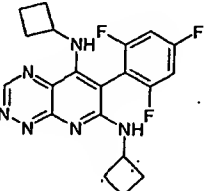
¹H NMR (CDCl₃) δ ppm: 1.20 (d,3H), 2.80 (s,3H), 4.10 (m,1H), 4.90 (bd,1H), 7.05 (m,1H), 7.60 (m,2H)

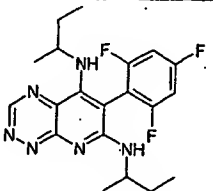
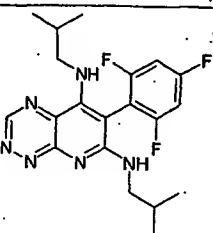
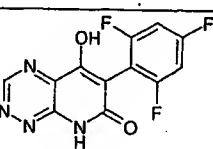
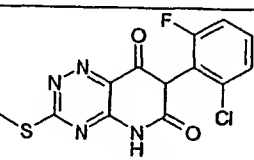
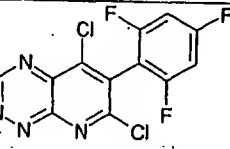
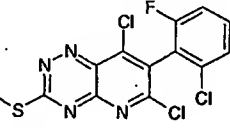
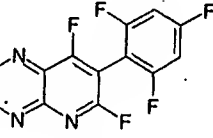
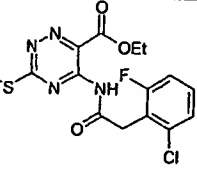
5 [8-chloro-3-methylsulfanyl-7-(2-chloro-6-fluorophenyl)-pyrido[2,3-e][1,2,4]triazin-6-yl]-isopropylamine

¹H NMR (CDCl₃) δ ppm: 1.35 (d,6H), 2.80 (s,3H), 4.70 (m,1H), 7.25 (m,2H), 7.45 (m,2H).

Table 136

| Compound No. | Table No. | Compound Structure | NMR data (ppm, in CDCl ₃ , unless otherwise stated) or Mpt. |
|--------------|-----------|--------------------|------------------------------------------------------------------------|
| 3 | 11 | | 1.15 (d,6H), 3.5 (b,1H), 6.7 (bflat,1H), 6.9 (t,2H), 9.85 (s,1H). |
| 3 | 16 | | 1.3 (d,6H), 4.75 (m,1H), 4.85 (bd,1H), 7.0 (t,2H), 9.80 (s,1H). |
| 3 | 26 | | 92-93°C |
| 5 | 26 | | 116-117°C |
| 12 | 26 | | 94°C |
| 23 | 26 | | 104°C |

| | | | |
|------|-----|-------------------------------------------------------------------------------------|-------------------------------------------------------------------------------|
| 58 | 26 |  | 129°C |
| 161 | 26 |  | 136-138°C |
| 2651 | 112 |  | 1.20 (d,3H), 2.80 (s,3H), 4.10 (m,1H), 4.90 (bd,1H), 7.05 (m,1H), 7.60 (m,2H) |
| 5961 | 112 |  | 1.35 (d,6H), 2.80 (s,3H), 4.70 (m,1H), 7.25 (m,2H), 7.45 (m,2H). |
| 161 | 122 |  | 190-192°C |
| 3 | 129 |  | 166-167°C |
| 14 | 129 |  | >220°C |
| 15 | 129 |  | >220°C |

| | | | |
|----|-----|-------------------------------------------------------------------------------------|------------------------------------------------------------------------------------------------|
| 23 | 129 |  | 0.65 -1.6 (m,16H), 2.85 (m,1H), 4.3 (bd,1H), 4.6 (m,1H), 6.25 (bd,1H), 6.90 (m,2H) 9.30 (s,1H) |
| 58 | 129 |  | 178°C |
| 13 | 132 |  | 6.95(t, 2H), 9.65(s, 1H). |
| 25 | 132 |  | (D ₆ -DMSO): 3.5 (s,3H), 7.0 (m,1H), 7.2 (m,2H). |
| 13 | 133 |  | 143-144°C |
| 25 | 133 |  | 2.8 (s,3H), 7.35 (m,1H), 7.45 (d,1H), 7.55 (m,1H). |
| 13 | 134 |  | 160-161°C |
| 25 | 135 |  | 1.45 (t,3H), 2.65 (s,3H), 4.5 (q,2H), 7.0 (m,1H), 7.2 (m,2H). |

EXAMPLE 4

This Example illustrates the fungicidal properties of the compounds of the general formula (1).

Compounds were tested in a leaf disk assay, with methods described below. Test compounds were dissolved in DMSO, and diluted into water to 200 ppm.

Erysiphe graminis f.sp. hordei (barley powdery mildew): barley leaf segments were placed on agar in a 24-well plate and sprayed with a solution of the test compound. After allowing to dry completely, for between 12 and 24 hours, the leaf disks were inoculated with a spore suspension of the fungus. After appropriate incubation the activity of a compound was assessed four days after inoculation as preventive fungicidal activity.

Erysiphe graminis f.sp. tritici (wheat powdery mildew): wheat leaf segments were placed on agar in a 24-well plate and sprayed with a solution of the test compound. After allowing to dry completely, for between 12 and 24 hours, the leaf disks were inoculated with a spore suspension of the fungus. After appropriate incubation the activity of a compound was assessed four days after inoculation as preventive fungicidal activity.

Puccinia recondita f.sp. tritici (wheat brown rust): wheat leaf segments were placed on agar in a 24-well plate and sprayed with a solution of the test compound. After allowing to dry completely, for between 12 and 24 hours, the leaf disks were inoculated with a spore suspension of the fungus. After appropriate incubation the activity of a compound was assessed nine days after inoculation as preventive fungicidal activity.

Pyrenophora teres (barley net blotch): barley leaf segments were placed on agar in a 24-well plate and sprayed with a solution of the test compound. After allowing to dry completely, for between 12 and 24 hours, the leaf disks were inoculated with a spore suspension of the fungus. After appropriate incubation the activity of a compound was assessed four days after inoculation as preventive fungicidal activity.

Pyricularia oryzae (rice blast): rice leaf segments were placed on agar in a 24-well plate and sprayed with a solution of the test compound. After allowing to dry completely, for between 12 and 24 hours, the leaf disks were inoculated with a spore suspension of the fungus. After appropriate incubation the activity of a compound was assessed four days after inoculation as preventive fungicidal activity.

The following compounds gave greater than 60% control of disease (number of compound first, followed by table number in brackets):

Erysiphe graminis f.sp. hordei, Compounds 3 (26), 23 (26), 58 (26), 23 (129), 58 (129);

Erysiphe graminis f.sp. tritici, Compounds 3 (11), 3 (16), 3 (26), 12 (26), 23 (26), 58 (26);

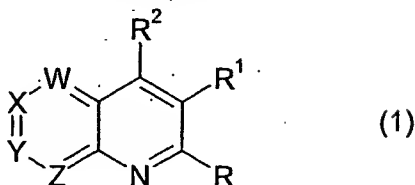
5 *Puccinia recondita f.sp. tritici*, Compounds 3 (11), 3 (16),

Pyrenophora teres, Compounds 12 (26), 23 (26), 58 (26);

Pyricularia oryzae, Compounds 3 (26), 12 (26), 23 (26), 58 (26).

CLAIMS

1. The compound of the general formula (1):



5 wherein

W, Z and one of X and Y are N and the other one of X and Y is CR⁸;

R⁸ is H, halo, C₁₋₄ alkyl, C₁₋₄ alkoxy, C₁₋₄ alkylthio or halo(C₁₋₄)alkyl;

R and R² are independently H, halo, C₁₋₈ alkyl, C₁₋₈ alkoxy, C₁₋₈ alkylthio, C₂₋₈ alkenyl, C₂₋₈ alkynyl, cyano or NR³R⁴, provided that at least one of R and R² is NR³R⁴;

10 R¹ is halo, C₁₋₈ alkyl, C₂₋₈ alkenyl, C₂₋₈ alkynyl, C₃₋₈ cycloalkyl, C₃₋₈ cycloalkyl(C₁₋₆)alkyl, C₁₋₈ alkoxy, C₁₋₈ alkylthio, aryl, aryloxy, arylthio, heteroaryl, heteroaryloxy, heteroarylthio, aryl(C₁₋₄)alkyl, aryl(C₁₋₄)alkoxy, heteroaryl(C₁₋₄)alkyl, heteroaryl(C₁₋₄)alkoxy, aryl(C₁₋₄)alkylthio, heteroaryl(C₁₋₄)alkylthio, morpholino, piperidino or pyrrolidino;

R³ and R⁴ are independently H, C₁₋₈ alkyl, C₂₋₈ alkenyl, C₂₋₈ alkynyl, aryl, aryl(C₁₋₈)alkyl, C₃₋₈ cycloalkyl, C₃₋₈ cycloalkyl(C₁₋₆)alkyl, heteroaryl, heteroaryl(C₁₋₈)alkyl, NR⁵R⁶, provided that
15 not both R³ and R⁴ are H or NR⁵R⁶, or

R³ and R⁴ together form a C₃₋₇ alkylene or C₃₋₇ alkenylene chain optionally substituted with one or more C₁₋₄ alkyl or C₁₋₄ alkoxy groups, or,

together with the nitrogen atom to which they are attached, R³ and R⁴ form a morpholine,

20 thiomorpholine, thiomorpholine S-oxide or thiomorpholine S-dioxide ring or a piperazine or piperazine N-(C₁₋₄)alkyl (especially N-methyl) ring; and

R⁵ and R⁶ are independently H, C₁₋₈ alkyl, C₂₋₈ alkenyl, C₂₋₈ alkynyl, aryl, aryl(C₁₋₈)alkyl, C₃₋₈ cycloalkyl, C₃₋₈ cycloalkyl(C₁₋₆)alkyl, heteroaryl or heteroaryl(C₁₋₈)alkyl;

any of the foregoing alkyl, alkenyl, alkynyl or cycloalkyl groups or moieties (other than for

25 R⁸) being optionally substituted with halogen, cyano, C₁₋₆ alkoxy, C₁₋₆ alkylcarbonyl, C₁₋₆ alkoxy carbonyl, C₁₋₆ haloalkoxy, C₁₋₆ alkylthio, tri(C₁₋₄)alkylsilyl, C₁₋₆ alkylamino or C₁₋₆ dialkylamino,

any of the foregoing morpholine, thiomorpholine, piperidine, piperazine and pyrrolidine rings being optionally substituted with C₁₋₄ alkyl (especially methyl), and

- any of the foregoing aryl or heteroaryl groups or moieties being optionally substituted with one or more substituents selected from halo, hydroxy, mercapto, C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₁₋₆ alkoxy, C₂₋₆ alkenyloxy, C₂₋₆ alkynyloxy, halo(C₁₋₆)alkyl, halo(C₁₋₆)alkoxy, C₁₋₆ alkylthio, halo(C₁₋₆)alkylthio, hydroxy(C₁₋₆)alkyl, C₁₋₄ alkoxy(C₁₋₆)alkyl, C₃₋₆ cycloalkyl, C₃₋₆ cycloalkyl(C₁₋₄)alkyl, phenoxy, benzyloxy, benzoyloxy, cyano, isocyano, thiocyanato, isothiocyanato, nitro, -NR^{'''}R^{'''}, -NHCOR^{'''}, -NHCONR^{'''}R^{'''}, -CONR^{'''}R^{'''}, -SO₂R^{'''}, -OSO₂R^{'''}, -COR^{'''}, -CR^{'''}=NR^{'''} or -N=CR^{'''}R^{'''}, in which R^{'''} and R^{'''} are independently hydrogen, C₁₋₄ alkyl, halo(C₁₋₄)alkyl, C₁₋₄ alkoxy, halo(C₁₋₄)alkoxy, C₁₋₄ alkylthio, C₃₋₆ cycloalkyl, C₃₋₆ cycloalkyl(C₁₋₄)alkyl, phenyl or benzyl, the phenyl and benzyl groups being optionally substituted with halogen, C₁₋₄ alkyl or C₁₋₄ alkoxy.
2. A compound according to claim 1 wherein W, Z and one of X and Y are N and the other one of X and Y is CH.
 3. A compound according to claim 1 or 2 wherein R² is NR³R⁴.
 4. A compound according to claim 3 wherein R is halo.
 5. A compound according to any one of the preceding claims wherein R³ is C₁₋₈ alkyl, halo(C₁₋₈)alkyl, hydroxy(C₁₋₈)alkyl, C₁₋₄ alkoxy(C₁₋₈)alkyl, C₁₋₄ alkoxyhalo(C₁₋₈)alkyl, tri(C₁₋₄)alkylsilyl(C₁₋₆)alkyl, C₁₋₄ alkylcarbonyl(C₁₋₈)alkyl, C₁₋₄ alkylcarbonylhalo(C₁₋₈)alkyl, phenyl(C₁₋₄)alkyl, C₂₋₈ alkenyl, halo(C₂₋₈)alkenyl, C₂₋₈ alkynyl, C₃₋₈ cycloalkyl optionally substituted with chloro, fluoro or methyl, C₃₋₈ cycloalkyl(C₁₋₄)alkyl, phenylamino, piperidino or morpholino, the phenyl ring of phenylalkyl or phenylamino being optionally substituted with one, two or three substituents selected from halo, C₁₋₄ alkyl, halo(C₁₋₄)alkyl, C₁₋₄ alkoxy and halo(C₁₋₄)alkoxy; and R⁴ is H, C₁₋₄ alkyl, halo(C₁₋₄)alkyl or amino, or R³ and R⁴ together form a C₃₋₇ alkylene or alkenylene chain optionally substituted with methyl, or, together with the nitrogen atom to which they are attached, R³ and R⁴ form a

morpholine, thiomorpholine, thiomorpholine S-oxide or thiomorpholine S-dioxide ring or a piperazine or piperazine *N*-(C₁₋₄)alkyl (especially *N*-methyl) ring, in which the morpholine or piperazine rings are optionally substituted with methyl.

- 5 6. A compound according to any one of the preceding claims wherein
 R^1 is phenyl optionally substituted with from one to five halogen atoms or with from one to three substituents selected from halo, C₁₋₄ alkyl, halo(C₁₋₄)alkyl, C₁₋₄ alkoxy or halo(C₁₋₄)alkoxy, pyridyl optionally substituted with from one to four halogen atoms or with from one to three substituents selected from halo, C₁₋₄ alkyl, halo(C₁₋₄)alkyl, C₁₋₄ alkoxy or halo(C₁₋₄)alkoxy, 2- or 3-thienyl optionally substituted with from one to three halogen atoms or with from one to three substituents selected from halo, C₁₋₄ alkyl, halo(C₁₋₄)alkyl, C₁₋₄ alkoxy or halo(C₁₋₄)alkoxy, or piperidino or morpholino both optionally substituted with one or two methyl groups.
- 10 7. A compound according to claim 6 wherein R^1 is 2,6-difluorophenyl, 2-fluoro-6-chlorophenyl, 2,5,6-trifluorophenyl, 2,4,6-trifluorophenyl, 2,6-difluoro-4-methoxyphenyl or pentafluorophenyl.
- 15 8. A compound according to claim 1 wherein
 20 W, Z and one of X and Y are N and the other one of X and Y is CR⁸;
 R^8 is H, halo, C₁₋₄ alkyl, C₁₋₄ alkoxy, C₁₋₄ alkylthio or halo(C₁₋₄)alkyl;
 one of R and R² (preferably R²) is NR³R⁴ and the other is halo;
 R^1 is halo, C₁₋₈ alkyl, C₂₋₈ alkenyl, C₂₋₈ alkynyl, C₃₋₈ cycloalkyl, C₃₋₈ cycloalkyl(C₁₋₆)-alkyl, C₁₋₈ alkoxy, C₁₋₈ alkylthio, aryl, aryloxy, arylthio, heteroaryl, heteroaryloxy,
 25 heteroarylthio, aryl(C₁₋₄)alkyl, aryl(C₁₋₄)alkoxy, heteroaryl(C₁₋₄)alkyl, heteroaryl-(C₁₋₄)alkoxy, aryl(C₁₋₄)alkylthio, heteroaryl(C₁₋₄)alkylthio, morpholino, piperidino or pyrrolidino;
 R^3 and R^4 are independently H, C₁₋₈ alkyl, C₂₋₈ alkenyl, C₂₋₈ alkynyl, aryl, aryl(C₁₋₈)-alkyl, C₃₋₈ cycloalkyl, C₃₋₈ cycloalkyl(C₁₋₆)alkyl, heteroaryl, heteroaryl(C₁₋₈)alkyl,
 30 NR⁵R⁶, provided that not both R^3 and R^4 are H or NR⁵R⁶, or
 R^3 and R^4 together form a C₃₋₇ alkylene or C₃₋₇ alkenylene chain optionally substituted with one or more C₁₋₄ alkyl or C₁₋₄ alkoxy groups, or,

together with the nitrogen atom to which they are attached, R^3 and R^4 form a morpholine, thiomorpholine, thiomorpholine S-oxide or thiomorpholine S-dioxide ring or a piperazine or piperazine N -(C_{1-4})alkyl (especially N -methyl) ring; and R^5 and R^6 are independently H, C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, aryl, aryl(C_{1-8})-alkyl, C_{3-8} cycloalkyl, C_{3-8} cycloalkyl(C_{1-6})alkyl, heteroaryl or heteroaryl(C_{1-8})alkyl; any of the foregoing alkyl, alkenyl, alkynyl or cycloalkyl groups or moieties (other than for R^8) being optionally substituted with halogen, cyano, C_{1-6} alkoxy, C_{1-6} alkylcarbonyl, C_{1-6} alkoxycarbonyl, C_{1-6} haloalkoxy, C_{1-6} alkylthio, tri(C_{1-4})alkylsilyl, C_{1-6} alkylamino or C_{1-6} dialkylamino,

any of the foregoing morpholine, thiomorpholine, piperidine, piperazine and pyrrolidine rings being optionally substituted with C_{1-4} alkyl (especially methyl), and any of the aryl, heteroaryl, aryloxy or heteroaryl groups being optionally substituted with one or more substituents selected from halo, hydroxy, mercapto, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{1-6} alkoxy, C_{2-6} alkenyloxy, C_{2-6} alkynyloxy, halo(C_{1-6})alkyl, halo(C_{1-6})alkoxy, C_{1-6} alkylthio, halo(C_{1-6})alkylthio, hydroxy(C_{1-6})alkyl, C_{1-4} alkoxy(C_{1-6})alkyl, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl(C_{1-4})alkyl, phenoxy, benzyloxy, benzoyloxy, cyano, isocyano, thiocyanato, isothiocyanato, nitro, $-NR^mR^n$, $-NHCOR^m$, $-NHCONR^mR^n$, $-CONR^mR^n$, $-SO_2R^m$, $-OSO_2R^m$, $-COR^m$, $-CR^m=NR^n$ or $-N=CR^mR^n$, in which R^m and R^n are independently hydrogen, C_{1-4} alkyl, halo(C_{1-4})alkyl, C_{1-4} alkoxy, halo(C_{1-4})alkoxy, C_{1-4} alkylthio, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl(C_{1-4})alkyl, phenyl or benzyl, the phenyl and benzyl groups being optionally substituted with halogen, C_{1-4} alkyl or C_{1-4} alkoxy.

9. A compound according to claim 1 wherein

W, Z and one of X and Y are N and the other one of X and Y is CR^8 ;
 R^8 is H, halo, C_{1-4} alkyl, C_{1-4} alkoxy, C_{1-4} alkylthio or halo(C_{1-4})alkyl;
 one of R and R^2 (preferably R^2) is NR^3R^4 and the other is halo;
 R^1 is halo, C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{3-8} cycloalkyl, C_{3-8} cycloalkyl(C_{1-6})-alkyl, C_{1-8} alkoxy, C_{1-8} alkylthio, aryl, aryloxy, arylthio, heteroaryl, heteroaryloxy, heteroarylthio, aryl(C_{1-4})alkyl, aryl(C_{1-4})alkoxy, heteroaryl(C_{1-4})alkyl, heteroaryl(C_{1-4})alkoxy, aryl(C_{1-4})alkylthio, heteroaryl(C_{1-4})alkylthio, morpholino, piperidino or pyrrolidino;

R^3 is C_{1-4} alkyl, halo(C_{1-4})alkyl, C_{2-4} alkenyl, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl(C_{1-4})-alkyl or phenylamino in which the phenyl ring is optionally substituted with one, two or three substituents selected from halo, C_{1-4} alkyl, halo(C_{1-4})alkyl, C_{1-4} alkoxy and halo(C_{1-4})alkoxy; and R^4 is H, C_{1-4} alkyl or amino, or

5 R^3 and R^4 together form a C_{4-6} alkylene chain optionally substituted with C_{1-4} alkyl or C_{1-4} alkoxy, or,

together with the nitrogen atom to which they are attached, R^3 and R^4 form a morpholine, thiomorpholine, thiomorpholine S-oxide or thiomorpholine S-dioxide ring or a piperazine or piperazine *N*-(C_{1-4})alkyl (especially *N*-methyl) ring;

10 any of the alkyl, alkenyl, alkynyl or cycloalkyl groups or moieties (other than for R^8) being optionally substituted with halogen, cyano, C_{1-6} alkoxy, C_{1-6} alkylcarbonyl, C_{1-6} alkoxy carbonyl, C_{1-6} haloalkoxy, C_{1-6} alkylthio, tri(C_{1-4})alkylsilyl, C_{1-6} alkylamino or C_{1-6} dialkylamino,

any of the foregoing morpholine, thiomorpholine, piperidine, piperazine and
15 pyrrolidine rings being optionally substituted with C_{1-4} alkyl (especially methyl), and any of the aryl or heteroaryl groups or moieties being optionally substituted with one or more substituents selected from halo, hydroxy, mercapto, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{1-6} alkoxy, C_{2-6} alkenyloxy, C_{2-6} alkynyloxy, halo(C_{1-6})alkyl, halo(C_{1-6})alkoxy, C_{1-6} alkylthio, halo(C_{1-6})alkylthio, hydroxy(C_{1-6})alkyl, C_{1-4} alkoxy(C_{1-6})-
20 alkyl, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl(C_{1-4})alkyl, phenoxy, benzyloxy, benzoyloxy, cyano, isocyano, thiocyanato, isothiocyanato, nitro, $-NR''''$, $-NHCOR''''$, $-NHCONR''''$, $-CONR''''$, $-SO_2R''''$, $-OSO_2R''''$, $-COR''''$, $-CR''''=NR''''$ or $-N=CR''''$, in which R'''' and R'''' are independently hydrogen, C_{1-4} alkyl, halo(C_{1-4})alkyl, C_{1-4} alkoxy, halo(C_{1-4})alkoxy, C_{1-4} alkylthio, C_{3-6} cycloalkyl, C_{3-6}
25 cycloalkyl(C_{1-4})alkyl, phenyl or benzyl, the phenyl and benzyl groups being optionally substituted with halogen, C_{1-4} alkyl or C_{1-4} alkoxy.

10. A compound according to claim 1 wherein

W, Z and one of X and Y are N and the other one of X and Y is CR^8 ;

30 R^8 is H, halo, C_{1-4} alkyl, C_{1-4} alkoxy, C_{1-4} alkylthio or halo(C_{1-4})alkyl;

R and R^2 are independently H, halo, C_{1-8} alkyl, C_{1-8} alkoxy, C_{1-8} alkylthio, C_{2-8} alkenyl, C_{2-8} alkynyl, cyano or NR^3R^4 , provided that at least one of R and R^2

(preferably R^2) is NR^3R^4 ;

R^1 is optionally substituted phenyl;

R^3 and R^4 are independently H, C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, aryl, aryl(C_{1-8})-alkyl, C_{3-8} cycloalkyl, C_{3-8} cycloalkyl(C_{1-6})alkyl, heteroaryl, heteroaryl(C_{1-8})alkyl, NR^5R^6 , provided that not both R^3 and R^4 are H or NR^5R^6 , or

R^3 and R^4 together form a C_{3-7} alkylene or C_{3-7} alkenylene chain optionally substituted with one or more C_{1-4} alkyl or C_{1-4} alkoxy groups, or, together with the nitrogen atom to which they are attached, R^3 and R^4 form a morpholine, thiomorpholine, thiomorpholine S-oxide or thiomorpholine S-dioxide ring or a piperazine or piperazine *N*-(C_{1-4})alkyl (especially *N*-methyl) ring; and R^5 and R^6 are independently H, C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, aryl, aryl(C_{1-8})-alkyl, C_{3-8} cycloalkyl, C_{3-8} cycloalkyl(C_{1-6})alkyl, heteroaryl or heteroaryl(C_{1-8})alkyl; any of the alkyl, alkenyl, alkynyl or cycloalkyl groups or moieties (other than for R^8) being optionally substituted with halogen, cyano, C_{1-6} alkoxy, C_{1-6} alkylcarbonyl, C_{1-6} alkoxy carbonyl, C_{1-6} haloalkoxy, C_{1-6} alkylthio, tri(C_{1-4})alkylsilyl, C_{1-6} alkylamino or C_{1-6} dialkylamino,

any of the foregoing morpholine, thiomorpholine, piperidine, piperazine and pyrrolidine rings being optionally substituted with C_{1-4} alkyl (especially methyl), and any of the aryl or heteroaryl groups or moieties, including the phenyl group of R^1 ,

being optionally substituted with one or more substituents selected from halo, hydroxy, mercapto, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{1-6} alkoxy, C_{2-6} alkenyloxy, C_{2-6} alkynyloxy, halo(C_{1-6})alkyl, halo(C_{1-6})alkoxy, C_{1-6} alkylthio, halo(C_{1-6})alkylthio, hydroxy(C_{1-6})alkyl, C_{1-4} alkoxy(C_{1-6})alkyl, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl(C_{1-4})alkyl, phenoxy, benzyloxy, benzoyloxy, cyano, isocyano, thiocyanato, isothiocyanato, nitro, $-NR''R'''$, $-NHCOR''$, $-NHCONR''R'''$, $-CONR''R'''$, $-SO_2R''$, $-OSO_2R''$, $-COR''$, $-CR''=NR'''$ or $-N=CR''R'''$, in which R'' and R''' are independently hydrogen, C_{1-4} alkyl, halo(C_{1-4})alkyl, C_{1-4} alkoxy, halo(C_{1-4})alkoxy, C_{1-4} alkylthio, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl(C_{1-4})alkyl, phenyl or benzyl, the phenyl and benzyl groups being optionally substituted with halogen, C_{1-4} alkyl or C_{1-4} alkoxy.

11. A compound according to claim 1 wherein

W, Z and one of X and Y are N and the other one of X and Y is CR^8 ;

R^8 is H, halo, C_{1-4} alkyl, C_{1-4} alkoxy, C_{1-4} alkylthio or halo(C_{1-4})alkyl;

R is H, halo, C_{1-4} alkyl, C_{1-4} alkoxy or cyano;

R^1 is phenyl optionally substituted with from one to five halogen atoms or with from one to three substituents selected from halo, C_{1-4} alkyl, halo(C_{1-4})alkyl, C_{1-4} alkoxy or halo(C_{1-4})alkoxy, pyridyl optionally substituted with from one to four halogen atoms or with from one to three substituents selected from halo, C_{1-4} alkyl, halo(C_{1-4})alkyl, C_{1-4} alkoxy or halo(C_{1-4})alkoxy, 2- or 3-thienyl optionally substituted with from one to three halogen atoms or with from one to three substituents selected from halo, C_{1-4} alkyl, halo(C_{1-4})alkyl, C_{1-4} alkoxy or halo(C_{1-4})alkoxy, or piperidino or morpholino both optionally substituted with one or two methyl groups;

R^2 is NR^3R^4 ;

R^3 is C_{1-8} alkyl, halo(C_{1-8})alkyl, hydroxy(C_{1-8})alkyl, C_{1-4} alkoxy(C_{1-8})alkyl, C_{1-4} alkoxyhalo(C_{1-8})alkyl, tri(C_{1-4})alkylsilyl(C_{1-6})alkyl, C_{1-4} alkylcarbonyl(C_{1-8})alkyl, C_{1-4} alkylcarbonylhalo(C_{1-8})alkyl, phenyl($_{1-4}$)alkyl, C_{2-8} alkenyl, halo(C_{2-8})alkenyl, C_{2-8} alkynyl, C_{3-8} cycloalkyl optionally substituted with chloro, fluoro or methyl, C_{3-8} cycloalkyl(C_{1-4})alkyl, phenylamino, piperidino or morpholino, the phenyl ring of phenylalkyl or phenylamino being optionally substituted with one, two or three substituents selected from halo, C_{1-4} alkyl, halo(C_{1-4})alkyl, C_{1-4} alkoxy and halo(C_{1-4})alkoxy; and

R^4 is H, C_{1-4} alkyl, halo(C_{1-4})alkyl or amino, or

R^3 and R^4 together form a C_{3-7} alkylene or C_{3-7} alkenylene chain optionally substituted with methyl, or,

together with the nitrogen atom to which they are attached, R^3 and R^4 form a morpholine, thiomorpholine, thiomorpholine S-oxide or thiomorpholine S-dioxide ring or a piperazine or piperazine *N*-(C_{1-4})alkyl (especially *N*-methyl) ring, in which the morpholine or piperazine rings are optionally substituted with methyl

12. A compound according to claim 1 wherein

W, Z and one of X and Y are N and the other one of X and Y is CR^8 ;

R^8 is H, halo, C_{1-4} alkyl, C_{1-4} alkoxy or halo(C_{1-4})alkyl;

R is halo;

R^1 is phenyl optionally substituted with from one to five halogen atoms or with from

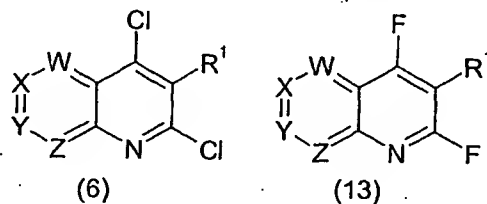
one to three substituents selected from halo, C₁₋₄ alkyl, halo(C₁₋₄)alkyl, C₁₋₄ alkoxy or halo(C₁₋₄)alkoxy;

R² is NR³R⁴;

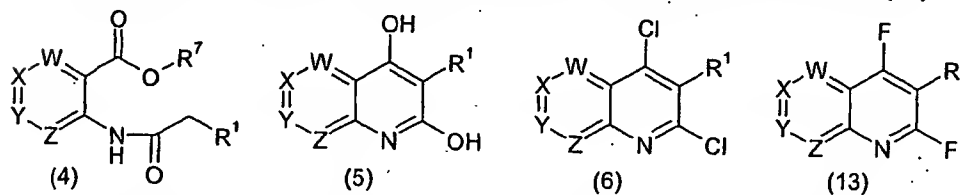
R³ is C₁₋₄ alkyl, halo(C₁₋₄)alkyl, C₂₋₄ alkenyl, C₃₋₆ cycloalkyl, C₃₋₆ cycloalkyl(C₁₋₄)-alkyl or phenylamino in which the phenyl ring is optionally substituted with one, two or three substituents selected from halo, C₁₋₄ alkyl, halo(C₁₋₄)alkyl, C₁₋₄ alkoxy and halo(C₁₋₄)alkoxy; and

R⁴ is H, C₁₋₄ alkyl or amino, or R³ and R⁴ together form a C₄₋₆ alkylene chain optionally substituted with methyl, or, together with the nitrogen atom to which they are attached, R³ and R⁴ form a morpholine ring.

13. A process for preparing a compound of the general formula (1) according to claim 1 wherein one of R and R² is chloro or fluoro and the other is NR³R⁴ and W, X, Y, Z, R¹, R³ and R⁴ are as defined in claim 1, which comprises reacting an amine of the general formula NR³R⁴ with a compound of the general formula (6) or (13):



14. The intermediate chemicals having the general formulae (4), (5), (6) and (13):



wherein W, X, Y, Z and R¹ are as defined in claim 1 and R⁷ is C₁₋₄ alkyl.

15. A plant fungicidal composition comprising a fungicidally effective amount of a compound as defined in claim 1 and a suitable carrier or diluent therefor.

16. A method of combating or controlling phytopathogenic fungi which comprises applying to a plant, to a seed of a plant, to the locus of the plant or seed or to soil or to any other plant growth medium, a fungicidally effective amount of a compound according to claim 1 or a composition according to claim 15.

INTERNATIONAL SEARCH REPORT

Application No

PCT/EP 03/05261

A. CLASSIFICATION OF SUBJECT MATTER

IPC 7 C07D487/04 A01N43/90

According to International Patent Classification (IPC) or to both national classification and IPC

B. FIELDS SEARCHED

Minimum documentation searched (classification system followed by classification symbols)

IPC 7 C07D A01N

Documentation searched other than minimum documentation to the extent that such documents are included in the fields searched

Electronic data base consulted during the international search (name of data base and, where practical, search terms used)

PAJ, WPI Data, EPO-Internal, CHEM ABS Data

C. DOCUMENTS CONSIDERED TO BE RELEVANT

| Category * | Citation of document, with indication, where appropriate, of the relevant passages | Relevant to claim No. |
|------------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|-----------------------|
| A | WO 02 088125 A (KRUEGER BERND-WIELAND ;MAULER-MACHNIK ASTRID (DE); DUNKEL RALF (DE) 7 November 2002 (2002-11-07) cited in the application page 5, line 2 -page 5, line 5; claims 1-7; examples 1-173 | 1-16 |
| A | WO 02 088126 A (KRUEGER BERND WIELAND ;MAULER-MACHNIK ASTRID (DE); DUNKEL RALF (DE) 7 November 2002 (2002-11-07) cited in the application page 5, line 6 -page 5, line 9; claims 1-7; examples 1-8323 | 1-16 |

☐ Further documents are listed in the continuation of box C.



Patent family members are listed in annex.

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Date of the actual completion of the international search

2 March 2004

Date of mailing of the international search report

17/03/2004

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INTERNATIONAL SEARCH REPORT

Application No
PCT/JP 03/05261

| Patent document cited in search report | | Publication date | Patent family member(s) | | Publication date |
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| WO 02088126 | A | 07-11-2002 | DE | 10121162 A1 | 31-10-2002 |
| | | | WO | 02088126 A1 | 07-11-2002 |

ABSTRACT

Acetamide derivatives represented by general formula (I) and physiologically acceptable acid-addition salts thereof, wherein X represents -O- or -NR₄; R₁ represents H, lower alkyl, lower alkenyl or cycloalkyl(lower)alkyl; R₂ represents lower alkyl, cycloalkyl, optionally substituted phenyl, etc.; R₃ represents H, lower alkyl or hydroxy(lower)alkyl; R₄ represents H, lower alkyl, etc.; R₅ represents H, lower alkyl, lower alkenyl, hydroxy(lower)alkyl, etc.; R₆ represents H, lower alkyl, CF₃ or optionally substituted phenyl, or R₅ and R₆ together form -(CH₂)_n-; R₇ represents H, halogeno, lower alkyl, lower alkoxy, CF₃, OH, NH₂, etc.; and R₈ represents H, halogeno, lower alkyl or lower alkoxy. The compounds act selectively on the peripheral BZ omega 3 receptor and exhibit excellent pharmacological effects, which makes them useful as a remedy and preventive for central diseases, for example, diseases associated with anxiety, depression and epilepsy, etc.

CLAIMS

No Claims were found.

DESCRIPTION

Text Not Available.